

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP093020\
 Data File : BP003415.D
 Acq On : 30 Sep 2020 16:02
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 01 10:05:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 10:05:32 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.534	152	101757	20.00	ng	0.00
21) Naphthalene-d8	10.310	136	386197	20.00	ng	# 0.00
39) Acenaphthene-d10	14.193	164	240014	20.00	ng	0.00
64) Phenanthrene-d10	16.945	188	540705	20.00	ng	# 0.00
76) Chrysene-d12	21.063	240	585519	20.00	ng	0.00
86) Perylene-d12	23.257	264	649108	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.152	112	478477	82.11	ng	0.00
7) Phenol-d6	6.740	99	639866	82.38	ng	0.00
23) Nitrobenzene-d5	8.687	82	543875	82.73	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	265935	72.64	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	1241300	75.64	ng	0.00
79) Terphenyl-d14	19.533	244	2074767	73.83	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.058	88	102811	42.93	ng	99
3) Pyridine	3.446	79	274392	43.03	ng	91
4) n-Nitrosodimethylamine	3.364	42	79300	41.58	ng	99
6) Aniline	6.869	93	360572	40.42	ng	94
8) 2-Chlorophenol	7.111	128	249109	38.38	ng	90
9) Benzaldehyde	6.681	77	165491	38.20	ng	87
10) Phenol	6.769	94	306917	41.37	ng	92
11) bis(2-Chloroethyl)ether	6.969	93	238368	40.54	ng	96
12) 1,3-Dichlorobenzene	7.428	146	295555	38.09	ng	# 95
13) 1,4-Dichlorobenzene	7.569	146	299271	38.13	ng	# 95
14) 1,2-Dichlorobenzene	7.887	146	280612	37.24	ng	97
15) Benzyl Alcohol	7.787	79	218254	42.28	ng	91
16) 2,2'-oxybis(1-Chloropr...	8.069	45	183471	42.92	ng	90
17) 2-Methylphenol	8.005	107	210646	39.26	ng	96
18) Hexachloroethane	8.605	117	114372	39.10	ng	93
19) n-Nitroso-di-n-propyla...	8.346	70	168704	40.54	ng	# 94
20) 3+4-Methylphenols	8.334	107	276793	38.77	ng	92
22) Acetophenone	8.358	105	365700	38.56	ng	95
24) Nitrobenzene	8.728	77	273750	41.62	ng	# 86
25) Isophorone	9.252	82	492054	40.84	ng	# 91
26) 2-Nitrophenol	9.440	139	137964	38.40	ng	# 82
27) 2,4-Dimethylphenol	9.522	122	194248	38.23	ng	90
28) bis(2-Chloroethoxy)met...	9.740	93	325091	40.54	ng	98
29) 2,4-Dichlorophenol	9.987	162	236625	37.27	ng	94
30) 1,2,4-Trichlorobenzene	10.187	180	274934	36.50	ng	99
31) Naphthalene	10.363	128	770715	37.78	ng	99
32) Benzoic acid	9.716	122	111986	34.08	ng	# 81
33) 4-Chloroaniline	10.481	127	334741	38.62	ng	# 93
34) Hexachlorobutadiene	10.657	225	181412	35.28	ng	99
35) Caprolactam	11.252	113	79385	37.47	ng	88
36) 4-Chloro-3-methylphenol	11.646	107	261820	38.25	ng	91
37) 2-Methylnaphthalene	11.987	142	544152	36.82	ng	97
38) 1-Methylnaphthalene	12.210	142	514119	36.64	ng	99
40) 1,2,4,5-Tetrachloroben...	12.375	216	290988	37.03	ng	99
41) Hexachlorocyclopentadiene	12.351	237	45724	27.48	ng	95
43) 2,4,6-Trichlorophenol	12.628	196	187117	36.83	ng	99

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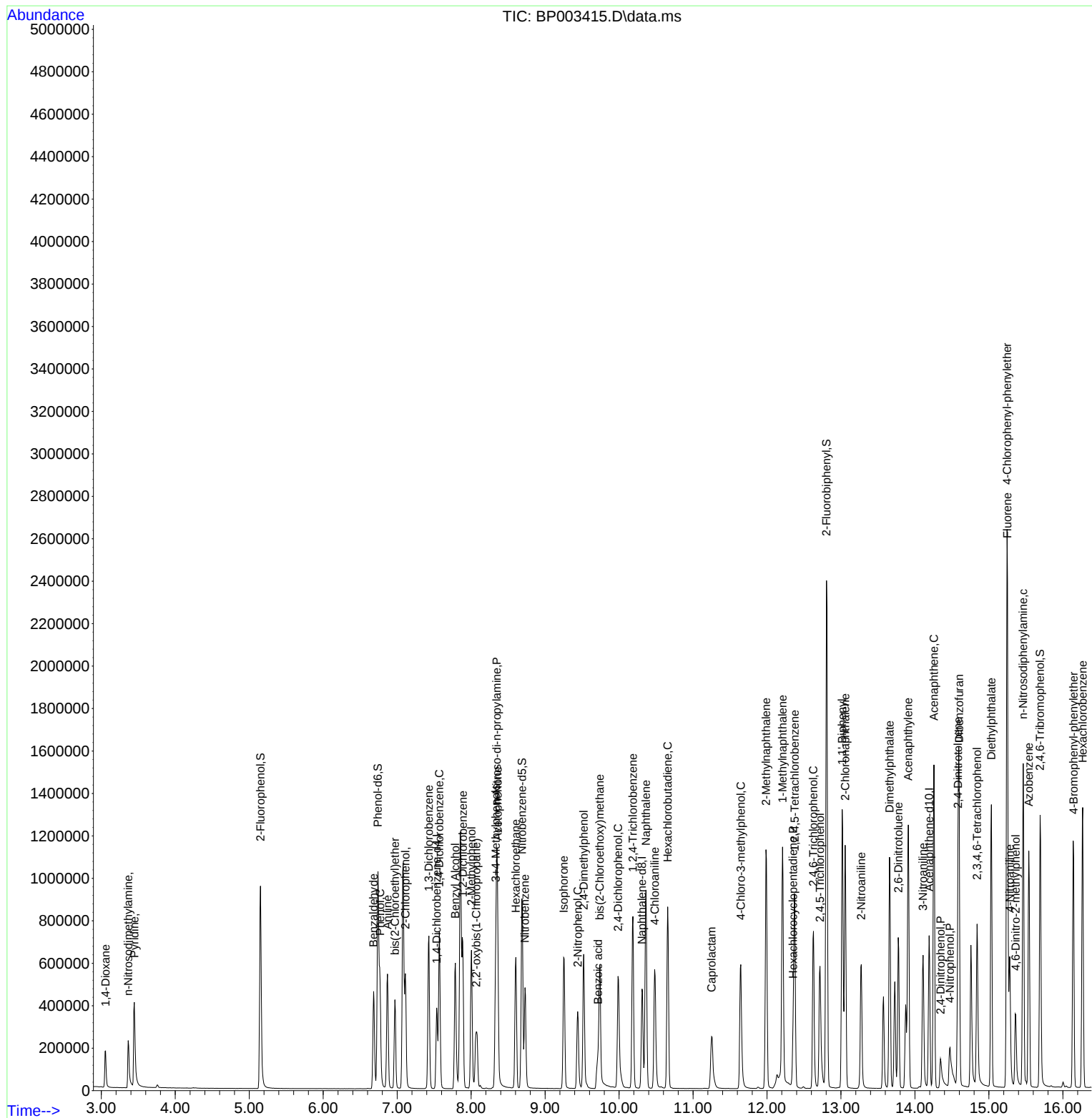
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	12.716	196	196408	37.58	ng	#	95
46) 1,1'-Biphenyl	13.016	154	689082	38.49	ng		98
47) 2-Chloronaphthalene	13.057	162	566722	38.23	ng		97
48) 2-Nitroaniline	13.275	65	164190	43.77	ng	#	76
49) Acenaphthylene	13.910	152	873265	38.56	ng		100
50) Dimethylphthalate	13.657	163	726877	37.93	ng		100
51) 2,6-Dinitrotoluene	13.775	165	157840	38.42	ng	#	76
52) Acenaphthene	14.257	154	525494	38.12	ng		100
53) 3-Nitroaniline	14.110	138	179806	40.39	ng	#	76
54) 2,4-Dinitrophenol	14.345	184	53497	32.28	ng	#	87
55) Dibenzofuran	14.598	168	821394	37.34	ng		96
56) 4-Nitrophenol	14.475	139	92317	33.93	ng	#	72
57) 2,4-Dinitrotoluene	14.581	165	217832	38.23	ng	#	67
58) Fluorene	15.251	166	651668	37.50	ng		99
59) 2,3,4,6-Tetrachlorophenol	14.840	232	181989	37.03	ng		98
60) Diethylphthalate	15.034	149	746569	38.19	ng		100
61) 4-Chlorophenyl-phenyle...	15.245	204	345976	36.47	ng		98
62) 4-Nitroaniline	15.281	138	194936	41.26	ng	#	74
63) Azobenzene	15.540	77	635249	42.63	ng		91
65) 4,6-Dinitro-2-methylph...	15.363	198	103104	36.14	ng		96
66) n-Nitrosodiphenylamine	15.463	169	595171	37.47	ng		100
67) 4-Bromophenyl-phenylether	16.145	248	238771	36.53	ng		98
68) Hexachlorobenzene	16.269	284	280370	36.11	ng		97
69) Atrazine	16.428	200	225453	36.19	ng		98
70) Pentachlorophenol	16.622	266	79888	28.71	ng		95
71) Phenanthrene	16.992	178	1099409	37.37	ng		99
72) Anthracene	17.081	178	1077861	37.18	ng		100
73) Carbazole	17.363	167	1033609	37.53	ng		99
74) Di-n-butylphthalate	17.928	149	1307887	37.15	ng	#	99
75) Fluoranthene	18.981	202	1368275	36.99	ng		99
77) Benzidine	19.163	184	564537	31.09	ng		99
78) Pyrene	19.333	202	1391163	38.10	ng		99
80) Butylbenzylphthalate	20.216	149	636472	38.10	ng	#	87
81) Benzo(a)anthracene	21.045	228	1398582	36.95	ng		99
82) 3,3'-Dichlorobenzidine	20.980	252	543855	37.18	ng		98
83) Chrysene	21.098	228	1358856	37.37	ng		99
84) Bis(2-ethylhexyl)phtha...	20.980	149	862300	36.66	ng		98
85) Di-n-octyl phthalate	21.839	149	1590787	37.02	ng	#	98
87) Indeno(1,2,3-cd)pyrene	25.498	276	1882905	38.27	ng		98
88) Benzo(b)fluoranthene	22.592	252	1517561	37.02	ng		99
89) Benzo(k)fluoranthene	22.645	252	1504456	38.30	ng		98
90) Benzo(a)pyrene	23.163	252	1464243	37.88	ng		98
91) Dibenzo(a,h)anthracene	25.510	278	1543958	38.03	ng		97
92) Benzo(g,h,i)perylene	26.198	276	1551024	38.27	ng		98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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