

Data Path : \\74.0.250.170\TERASTORAGE\SVOASRV\HPCHEM1\BNA\_F\DATA\BF062919\  
 Data File : BF115318.D  
 Acq On : 29 Jun 2019 13:06  
 Operator : HP/JU  
 Sample : K3531-03  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 FK-GF-02-043-B

Quant Time: Jun 30 02:14:48 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF062119.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Jun 28 05:26:43 2019  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	191760	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	752777	20.00	ng	0.00
39) Acenaphthene-d10	9.62	164	371445	20.00	ng	0.00
64) Phenanthrene-d10	11.10	188	755339	20.00	ng	0.00
76) Chrysene-d12	13.71	240	666036	20.00	ng	0.00
87) Perylene-d12	15.08	264	689439	20.00	ng	0.00

## System Monitoring Compounds

5) 2-Fluorophenol	5.19	112	1103614	88.81	ng	0.00
7) Phenol-d6	6.24	99	1406393	93.25	ng	0.00
23) Nitrobenzene-d5	7.16	82	827801	67.65	ng	0.00
42) 2,4,6-Tribromophenol	10.41	330	404032	103.15	ng	0.00
45) 2-Fluorobiphenyl	8.95	172	1560323	71.35	ng	0.00
79) Terphenyl-d14	12.68	244	1851497	59.10	ng	0.00

## Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.15	88	127	0.022	ng	# 1
3) Pyridine	3.00	79	115	0.007	ng	# 1
6) Aniline	6.25	93	292	0.014	ng	# 1
8) 2-Chlorophenol	6.38	128	322	0.023	ng	# 1
9) Benzaldehyde	6.24	77	2010	0.204	ng	# 1
10) Phenol	6.25	94	5604	0.317	ng	89
11) bis(2-Chloroethyl)ether	6.37	93	1590	0.122	ng	# 1
15) Benzyl Alcohol	6.75	79	13518	1.231	ng	# 38
22) Acetophenone	7.00	105	346	0.020	ng	# 52
24) Nitrobenzene	7.16	77	2517	0.187	ng	# 33
28) bis(2-Chloroethoxy)methane	7.88	93	144	0.009	ng	# 1
31) Naphthalene	7.90	128	5194	0.141	ng	99
33) 4-Chloroaniline	7.90	127	632	0.037	ng	# 47
37) 2-Methylnaphthalene	8.59	142	2092	0.084	ng	88
38) 1-Methylnaphthalene	8.69	142	1314	0.055	ng	# 83
46) 1,1'-Biphenyl	9.05	154	3419	0.115	ng	93
47) 2-Chloronaphthalene	9.35	162	923	0.038	ng	# 1
48) 2-Nitroaniline	9.35	65	1577	0.224	ng	# 1
49) Acenaphthylene	9.48	152	7355	0.196	ng	98
50) Dimethylphthalate	9.35	163	179834	6.581	ng	99
51) 2,6-Dinitrotoluene	9.35	165	1818	0.288	ng	# 18
52) Acenaphthene	9.65	154	2402	0.102	ng	90
55) Dibenzofuran	9.82	168	2978	0.088	ng	# 87
56) 4-Nitrophenol	9.82	139	1100	0.178	ng	# 11
60) Diethylphthalate	10.05	149	939	0.034	ng	# 83
63) Azobenzene	10.32	77	1231	0.048	ng	# 1
66) n-Nitrosodiphenylamine	10.28	169	621	0.026	ng	# 71
71) Phenanthrene	11.12	178	48548	1.194	ng	96
72) Anthracene	11.17	178	15632	0.381	ng	96
73) Carbazole	11.32	167	5637	0.154	ng	97
74) Di-n-butylphthalate	11.67	149	4636	0.109	ng	# 91
75) Fluoranthene	12.30	202	113830	2.805	ng	96
77) Benzidine	12.42	184	881	0.030	ng	# 1
78) Pyrene	12.52	202	117130	2.288	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
80) Butylbenzylphthalate	13.16	149	1142	0.051	ng	# 77
81) Benzo(a)anthracene	13.70	228	63169	1.322	ng	95
82) 3,3'-Dichlorobenzidine	13.68	252	935	0.054	ng	# 43
83) Chrysene	13.74	228	60580	1.384	ng	96
84) Bis(2-ethylhexyl)phthalate	13.72	149	6444	0.218	ng	# 88
85) Di-n-octyl phthalate	14.34	149	812	0.016	ng	# 1
86) Indeno(1,2,3-cd)pyrene	16.33	276	31663	0.611	ng	99
90) Benzo(a)pyrene	15.02	252	54289	1.400	ng	# 95
91) Dibenzo(a,h)anthracene	16.35	278	8715	0.241	ng	# 62
92) Benzo(g,h,i)perylene	16.71	276	28756	0.785	ng	95

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

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