

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060716\
 Data File : BE091895.D
 Acq On : 8 Jun 2016 3:11
 Operator : UM/SJ
 Sample : PB91033BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB91033BS

Quant Time: Jun 08 14:38:48 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 08 14:06:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	22673	5.00	ng	0.00
8) Naphthalene-d8	10.55	136	97455	5.00	ng	0.00
15) Acenaphthene-d10	14.38	164	57354	5.00	ng	0.00
24) Phenanthrene-d10	17.12	188	145632	5.00	ng	0.00
31) Chrysene-d12	21.27	240	143342	5.00	ng	-0.03
40) Perylene-d12	23.71	264	187510	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.37	112	34156	7.06	ng	0.00
5) Phenol-d6	6.94	99	48340	8.07	ng	-0.02
9) Nitrobenzene-d5	8.91	82	31076	4.24	ng	-0.02
16) 2,4,6-Tribromophenol	15.87	330	16329	9.99	ng	-0.02
17) 2-Fluorobiphenyl	13.01	172	68889	3.96	ng	0.00
34) Terphenyl-d14	19.74	244	119752	-5.00	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.30	88	8894	2.96	ng	90
3) n-Nitrosodimethylamine	3.61	42	12909	3.40	ng	99
6) bis(2-Chloroethyl)ether	7.20	93	23025	3.55	ng	98
7) n-Nitroso-di-n-propylamine	8.54	70	19030	3.71	ng	# 93
10) Nitrobenzene	8.96	77	26253	3.93	ng	# 66
11) bis(2-Chloroethoxy)methane	9.95	93	38036	3.85	ng	# 39
12) Naphthalene	10.59	128	78224	3.72	ng	98
13) Hexachlorobutadiene	10.88	225	11584	3.70	ng	99
14) 2-Methylnaphthalene	12.20	142	55795	4.11	ng	99
18) Acenaphthylene	14.10	152	96515	3.72	ng	# 91
19) 2,6-Dinitrotoluene	13.96	165	15683	4.04	ng	96
20) Acenaphthene	14.44	154	64760	4.08	ng	92
21) 2,4-Dinitrotoluene	14.76	165	18114	4.08	ng	# 1
22) Fluorene	15.42	166	75969	3.92	ng	100
23) 4-Chlorophenyl-phenylether	15.42	204	36196	3.80	ng	97
25) 4-Bromophenyl-phenylether	16.31	248	21472	4.10	ng	96
26) Hexachlorobenzene	16.44	284	22397	4.18	ng	98
27) Pentachlorophenol	16.78	266	23911	8.03	ng	93
28) Phenanthrene	17.16	178	138239	4.10	ng	99
29) Anthracene	17.25	178	131499	4.38	ng	100
30) Fluoranthene	19.17	202	170891	4.69	ng	98
32) Benzidine	19.36	184	65969	5.69	ng	# 94
33) Pyrene	19.53	202	186511	5.16	ng	99
35) Benzo(a)anthracene	21.27	228	844261m	4.92	ng	
36) 3,3'-Dichlorobenzidine	21.21	252	49961	4.32	ng	# 99
37) Chrysene	21.30	228	232609m	1.59	ng	
38) Bis(2-ethylhexyl)phthalate	21.19	149	76455	2.95	ng	# 100
39) Indeno(1,2,3-cd)pyrene	26.27	276	205310	4.70	ng	# 75
41) Benzo(b)fluoranthene	22.97	252	192568	4.20	ng	98
42) Benzo(k)fluoranthene	23.02	252	200356	4.21	ng	95
43) Benzo(a)pyrene	23.59	252	186775	4.35	ng	# 95
44) Dibenzo(a,h)anthracene	26.29	278	175442	4.38	ng	# 93
45) Benzo(a,h,i)perylene	27.05	276	176456	4.20	ng	# 93

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

