

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101617\
 Data File : BE094420.D
 Acq On : 16 Oct 2017 16:53
 Operator : SJ/JU
 Sample : I5266-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-BN-WP

Quant Time: Oct 16 20:09:55 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 14:54:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1425	0.40	ng	-0.01
7) Naphthalene-d8	10.69	136	7092	0.40	ng	-0.02
13) Acenaphthene-d10	14.53	164	3940	0.40	ng	0.00
19) Phenanthrene-d10	17.24	188	6841	0.40	ng	-0.03
27) Chrysene-d12	21.43	240	9050	0.40	ng	0.00
34) Perylene-d12	23.96	264	8825	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.51	112	759806	157.78	ng	0.00
5) Phenol-d6	7.11	99	1170808	155.89	ng	-0.01
8) Nitrobenzene-d5	9.06	82	770920	125.09	ng	-0.02
11) 2-Methylnaphthalene-d10	0.00	152	0d	0.00	ng	
14) 2,4,6-Tribromophenol	16.03	330	271276	180.73	ng	0.00
15) 2-Fluorobiphenyl	13.15	172	1533840	109.06	ng	0.00
25) Fluoranthene-d10	0.00	212	0d	0.00	ng	
29) Terphenyl-d14	19.87	244	1794728	98.28	ng	0.00

Target Compounds					Qvalue	
6) bis(2-Chloroethyl)ether	7.33	93	679723	122.371	ng	99
9) Naphthalene	10.76	128	9622039	121.329	ng	99
10) Hexachlorobutadiene	11.03	225	216836	68.404	ng	98
12) 2-Methylnaphthalene	12.36	142	1239632	94.816	ng	97
16) Acenaphthylene	14.24	152	1495294	69.984	ng	99
17) Acenaphthene	14.58	154	884975	65.782	ng	97
18) Fluorene	15.58	166	1702259	98.113	ng	98
23) Phenanthrene	17.30	178	305199	11.334	ng	96
26) Fluoranthene	19.32	202	1531914	45.847	ng	99
28) Pyrene	19.68	202	1504421	44.059	ng	98
31) Chrysene	21.47	228	1931427	64.103	ng	# 65
33) Indeno(1,2,3-cd)pyrene	26.70	276	2877585	94.824	ng	96
35) Benzo(b)fluoranthene	23.21	252	3757569	125.527	ng	# 39
36) Benzo(k)fluoranthene	23.25	252	1038693	35.218	ng	# 38
37) Benzo(a)pyrene	23.85	252	807943	28.717	ng	# 31
38) Dibenzo(a,h)anthracene	26.70	278	930160	34.926	ng	# 42
39) Benzo(g,h,i)perylene	27.55	276	3135411	122.401	ng	# 52

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

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