

Data Path : Z:\HPCHEM1\BNA E\Data\BE032318\
 Data File : BE095824.D
 Acq On : 23 Mar 2018 14:47
 Operator : SJ/JU
 Sample : J1888-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-BN-WP

Quant Time: Mar 23 23:50:21 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1235	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	4383	0.40	ng	0.00
13) Acenaphthene-d10	15.00	164	2562	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	6012	0.40	ng	0.00
27) Chrysene-d12	21.80	240	7185	0.40	ng	0.00
34) Perylene-d12	24.44	264	6826	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.91	112	440995	108.32	ng	0.00
5) Phenol-d6	7.57	99	642745	114.81	ng	0.00
8) Nitrobenzene-d5	9.61	82	467553	90.29	ng	0.01
11) 2-Methylnaphthalene-d10	12.79	152	2	0.00	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	207668	140.46	ng	0.00
15) 2-Fluorobiphenyl	13.64	172	862021	79.83	ng	0.00
25) Fluoranthene-d10	19.68	212	64	0.00	ng	-0.01
29) Terphenyl-d14	20.25	244	1327212	86.98	ng	0.01

Target Compounds					Qvalue	
9) Naphthalene	11.31	128	3840034	76.707	ng	99
12) 2-Methylnaphthalene	12.87	142	870963	101.872	ng	94
16) Acenaphthylene	14.74	152	2075800	141.670	ng	98
17) Acenaphthene	15.07	154	767812	85.914	ng	95
18) Fluorene	16.03	166	442404	37.633	ng	# 78
23) Phenanthrene	17.75	178	1287204	71.523	ng	99
26) Fluoranthene	19.72	202	2173670	86.955	ng	# 96
28) Pyrene	20.07	202	1487920	50.840	ng	98
30) Benzo(a)anthracene	21.79	228	1550473	62.227	ng	98
31) Chrysene	21.84	228	755472	33.226	ng	97
32) Bis(2-ethylhexyl)phthalate	21.64	149	2377199	31.196	ng	100
33) Indeno(1,2,3-cd)pyrene	27.29	276	1687690	68.891	ng	# 91
35) Benzo(b)fluoranthene	23.63	252	1823396	80.751	ng	# 85
36) Benzo(k)fluoranthene	23.68	252	1314299	58.190	ng	95
37) Benzo(a)pyrene	24.34	252	2579701	121.875	ng	# 91
38) Dibenzo(a,h)anthracene	27.30	278	1005287	49.988	ng	# 92
39) Benzo(a,h,i)perylene	28.17	276	1168306	57.839	ng	# 89

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

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