

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020718\
 Data File : BE095549.D
 Acq On : 7 Feb 2018 12:03
 Operator : SJ/JU
 Sample : J1196-02DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 1050DL

Quant Time: Feb 08 01:44:55 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 20:27:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1154	0.40	ng	0.00
7) Naphthalene-d8	11.28	136	4555	0.40	ng	0.00
13) Acenaphthene-d10	15.00	164	2668	0.40	ng	0.00
19) Phenanthrene-d10	17.71	188	6499	0.40	ng	0.00
27) Chrysene-d12	21.79	240	6814	0.40	ng	0.00
34) Perylene-d12	24.43	264	6405	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.92	112	108	0.03	ng	0.00
5) Phenol-d6	7.55	99	176	0.04	ng	-0.01
8) Nitrobenzene-d5	9.62	82	208	0.04	ng	0.00
11) 2-Methylnaphthalene-d10	12.80	152	308	0.04	ng	0.00
14) 2,4,6-Tribromophenol	16.47	330	39	0.03	ng	0.00
15) 2-Fluorobiphenyl	13.65	172	523	0.05	ng	0.00
25) Fluoranthene-d10	19.69	212	3974	0.05	ng	0.00
29) Terphenyl-d14	20.25	244	928	0.06	ng	0.00

Target Compounds				Qvalue		
6) bis(2-Chloroethyl)ether	7.83	93	106	0.024	ng	82
9) Naphthalene	11.31	128	70990	1.342	ng	99
12) 2-Methylnaphthalene	12.88	142	276	0.031	ng	# 91
16) Acenaphthylene	14.73	152	9415	0.619	ng	99
17) Acenaphthene	15.06	154	7951	0.841	ng	94
18) Fluorene	16.04	166	6684	0.562	ng	# 79
20) 4-Bromophenyl-phenylether	16.91	248	211	0.054	ng	# 80
23) Phenanthrene	17.74	178	32915	1.649	ng	98
24) Anthracene	17.83	178	4790	0.257	ng	95
26) Fluoranthene	19.72	202	21609	0.858	ng	# 97
28) Pvrene	20.06	202	23735	0.830	ng	95
30) Benzo(a)anthracene	21.78	228	12259	0.540	ng	98
31) Chrysene	21.84	228	5643	0.252	ng	98
32) Bis(2-ethylhexyl)phthalate	21.65	149	4392	0.084	ng	# 98
33) Indeno(1,2,3-cd)pvrene	27.23	276	8470	0.392	ng	96
35) Benzo(b)fluoranthene	23.61	252	9435	0.432	ng	# 91
36) Benzo(k)fluoranthene	23.66	252	18740	0.835	ng	96
37) Benzo(a)pvrene	24.31	252	13322	0.655	ng	93
38) Dibenzo(a,h)anthracene	27.26	278	10803	0.607	ng	96
39) Benzo(g,h,i)perylene	28.10	276	9379	0.491	ng	# 93

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

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