

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032618\
 Data File : BE095853.D
 Acq On : 26 Mar 2018 00:00
 Operator : SJ/JU
 Sample : J1994-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C-MW-10-031918

Quant Time: Mar 27 06:50:54 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	1208	0.40	ng	0.00
7) Naphthalene-d8	11.26	136	4601	0.40	ng	0.00
13) Acenaphthene-d10	14.99	164	2635	0.40	ng	-0.01
19) Phenanthrene-d10	17.70	188	6412	0.40	ng	-0.01
27) Chrysene-d12	21.79	240	7687	0.40	ng	-0.01
34) Perylene-d12	24.41	264	7589	0.40	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.91	112	620	0.16	ng	0.00
5) Phenol-d6	7.56	99	570	0.10	ng	0.00
8) Nitrobenzene-d5	9.59	82	2589	0.48	ng	0.00
11) 2-Methylnaphthalene-d10	12.79	152	2859	0.37	ng	-0.01
14) 2,4,6-Tribromophenol	16.46	330	667	0.44	ng	-0.01
15) 2-Fluorobiphenyl	13.63	172	6356	0.57	ng	-0.01
25) Fluoranthene-d10	19.68	212	36768	0.39	ng	0.00
29) Terphenyl-d14	20.24	244	10164	0.62	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
22) Pentachlorophenol	17.36	266	152	0.236	ng	# 80
23) Phenanthrene	17.74	178	2225	0.116	ng	100
26) Fluoranthene	19.71	202	4114	0.154	ng	97
28) Pyrene	20.06	202	2947	0.094	ng	# 88
30) Benzo(a)anthracene	21.77	228	908	0.034	ng	# 73
31) Chrysene	21.83	228	1493	0.061	ng	# 83
32) Bis(2-ethylhexyl)phthalate	21.65	149	6173	0.076	ng	99
33) Indeno(1,2,3-cd)pyrene	27.21	276	919	0.035	ng	# 79
35) Benzo(b)fluoranthene	23.61	252	1642	0.065	ng	# 64
39) Benzo(a,h,i)perylene	28.10	276	882	0.039	ng	# 51

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

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