

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102517\
 Data File : BE094534.D
 Acq On : 25 Oct 2017 15:46
 Operator : SJ/JU
 Sample : I5693-05DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-33(0.5-1.5)-100417DL

Quant Time: Oct 25 19:02:46 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 25 18:44:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1417	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	7241	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	5020	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9136	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	9435	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	9829	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	437	0.04	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	1977	0.07	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	39279	2.219	ng/ul#	95
5) 2-Methylnaphthalene	12.34	142	71689	5.788	ng/ul	99
7) Acenaphthylene	14.24	152	3481	0.153	ng/ul#	31
8) Acenaphthene	14.57	153	6056	0.371	ng/ul	92
9) Fluorene	15.56	166	7644	0.417	ng/ul#	84
11) Pentachlorophenol	17.03	266	29	0.064	ng/ul#	45
12) Phenanthrene	17.29	178	75688	3.137	ng/ul#	93
13) Anthracene	17.38	178	13471	0.566	ng/ul	96
15) Fluoranthene	19.31	202	79062	2.834	ng/ul	83
17) Pyrene	19.67	202	101612	3.484	ng/ul#	83
18) Benzo(a)anthracene	21.40	228	59017	2.188	ng/ul#	92
19) Chrysene	21.46	228	66062	2.370	ng/ul#	91
21) Benzo(b)fluoranthene	23.18	252	115398	4.174	ng/ul	89
22) Benzo(k)fluoranthene	23.18	252	114703	3.871	ng/ul#	91
23) Benzo(a)pyrene	23.84	252	60022	2.172	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.64	276	47783	1.671	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.64	278	12171	0.505	ng/ul#	63
26) Benzo(g,h,i)perylene	27.48	276	40359	1.758	ng/ul	91

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

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