

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042518\
 Data File : BE096142.D
 Acq On : 25 Apr 2018 10:25
 Operator : SJ/JU
 Sample : J2467-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3XS5

Quant Time: Apr 26 04:21:54 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 26 04:18:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	740	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2736	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1504	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4392	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	6755	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	5234	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1762	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	7253	0.36	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.23	128	1944	0.255	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	1434	0.281	ng/ul	100
7) Acenaphthylene	14.67	152	6524	0.841	ng/ul	97
8) Acenaphthene	15.01	153	1957	0.350	ng/ul	94
9) Fluorene	15.98	166	2570	0.347	ng/ul	92
11) Pentachlorophenol	17.32	266	237	0.224	ng/ul	96
12) Phenanthrene	17.70	178	9619	0.744	ng/ul#	89
13) Anthracene	17.79	178	3260	0.251	ng/ul#	93
15) Fluoranthene	19.67	202	8420	0.426	ng/ul	85
17) Pyrene	20.02	202	5320	0.393	ng/ul#	78
18) Benzo(a)anthracene	21.73	228	11457	0.560	ng/ul#	85
19) Chrysene	21.79	228	7684	0.409	ng/ul	96
21) Benzo(b)fluoranthene	23.54	252	6595	0.333	ng/ul	87
23) Benzo(a)pyrene	24.23	252	8561	0.497	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.14	276	7595	0.327	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.15	278	6737	0.347	ng/ul#	89
26) Benzo(g,h,i)perylene	27.99	276	7642	0.378	ng/ul	95

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

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