

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042418\
 Data File : BE096137.D
 Acq On : 25 Apr 2018 1:43
 Operator : SJ/JU
 Sample : J2467-10MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3XX6MS

Quant Time: Apr 25 06:57:52 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 25 05:27:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	816	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	3145	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1873	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	5401	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	7751	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	5947	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1869	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	7767	0.31	ng/ul	0.00
Target Compounds				Qvalue		
8) Acenaphthene	15.01	153	2055	0.295	ng/ul	96
9) Fluorene	15.98	166	188	0.020	ng/ul#	88
11) Pentachlorophenol	17.32	266	431	0.332	ng/ul	96
12) Phenanthrene	17.70	178	1028	0.065	ng/ul#	91
15) Fluoranthene	19.67	202	1181	0.049	ng/ul	86
17) Pyrene	20.02	202	4679	0.301	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	495	0.021	ng/ul#	62
19) Chrysene	21.79	228	562	0.026	ng/ul#	75
21) Benzo(b)fluoranthene	23.54	252	729	0.032	ng/ul#	60
24) Indeno(1,2,3-cd)pyrene	27.13	276	596	0.023	ng/ul#	85
26) Benzo(g,h,i)perylene	28.00	276	544	0.024	ng/ul#	52

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

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