

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070918\
 Data File : BE096888.D
 Acq On : 9 Jul 2018 19:24
 Operator : SJ/JU
 Sample : J3765-05DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H07DL

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:56:43 PM

Quant Time: Jul 10 03:51:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 10 02:37:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3894	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	21297	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	13142	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	30801	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	32786	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	32056m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	1074	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	2857	0.04	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.19	128	9938	0.193	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	3568	0.102	ng/ul	99
7) Acenaphthylene	13.74	152	41161	0.764	ng/ul	96
8) Acenaphthene	14.09	153	1574	0.035	ng/ul#	88
9) Fluorene	15.08	166	1802	0.039	ng/ul#	88
12) Phenanthrene	16.82	178	37823	0.476	ng/ul#	89
13) Anthracene	16.91	178	53980	0.800	ng/ul#	92
15) Fluoranthene	18.85	202	250587	2.962	ng/ul	84
17) Pyrene	19.21	202	349431	3.247	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	292008	3.776	ng/ul#	82
19) Chrysene	21.02	228	396306	3.956	ng/ul	98
21) Benzo(b)fluoranthene	22.55	252	1031059m	11.785	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	307533m	3.220	ng/ul	
23) Benzo(a)pyrene	23.12	252	506751	6.848	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.51	276	437347	5.693	ng/ul#	74
25) Dibenzo(a,h)anthracene	25.52	278	101956m	1.671	ng/ul	
26) Benzo(g,h,i)perylene	26.21	276	378098	5.071	ng/ul#	92

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

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