

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070918\
 Data File : BE096884.D
 Acq On : 9 Jul 2018 16:54
 Operator : SJ/JU
 Sample : J3721-19DL 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H05DL

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:56:33 PM

Quant Time: Jul 09 17:41:48 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 : Sat Jul 07 07:26:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3025	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	16333	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	10161	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	23975	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	22568	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	22295m	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	1809	0.07	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	4496	0.08	ng/ul	0.00

Target Compounds				Qvalue		
7) Acenaphthylene	13.74	152	3114	0.075	ng/ul	99
12) Phenanthrene	16.82	178	2365	0.038	ng/ul#	91
13) Anthracene	16.91	178	3454	0.066	ng/ul	92
15) Fluoranthene	18.84	202	34334	0.521	ng/ul	84
17) Pyrene	19.21	202	48781	0.658	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	22525	0.423	ng/ul#	83
19) Chrysene	21.02	228	39160	0.568	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	77134m	1.268	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	24850m	0.374	ng/ul	
23) Benzo(a)pyrene	23.12	252	34820	0.677	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	22152	0.415	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.52	278	5489	0.129	ng/ul	92
26) Benzo(g,h,i)perylene	26.21	276	17036	0.329	ng/ul	94

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

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