

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096867.D
 Acq On : 6 Jul 2018 23:42
 Operator : SJ/JU
 Sample : J3721-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H20

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:12 PM

Quant Time: Jul 09 05:36:27 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 : Fri Jul 06 16:55:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	3835	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	20405	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	12989	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	30993m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	30547m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	28318m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8484	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	23098m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	3459	0.070	ng/ul#	95
5) 2-Methylnaphthalene	11.81	142	961	0.029	ng/ul	98
7) Acenaphthylene	13.74	152	69109	1.298	ng/ul	96
8) Acenaphthene	14.09	153	2121	0.048	ng/ul#	86
9) Fluorene	15.08	166	6645	0.147	ng/ul	91
12) Phenanthrene	16.82	178	39905m	0.499	ng/ul	
13) Anthracene	16.91	178	75224m	1.109	ng/ul	
15) Fluoranthene	18.85	202	249912m	2.936	ng/ul	
17) Pyrene	19.21	202	693404	6.915	ng/ul#	82
18) Benzo(a)anthracene	20.97	228	632839m	8.783	ng/ul	
19) Chrysene	21.02	228	1023925	10.969	ng/ul	96
21) Benzo(b)fluoranthene	22.55	252	2340269m	30.280	ng/ul	
22) Benzo(k)fluoranthene	22.59	252	718052m	8.510	ng/ul	
23) Benzo(a)pyrene	23.13	252	1318999m	20.177	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.52	276	834949	12.304	ng/ul#	75
25) Dibenzo(a,h)anthracene	25.52	278	214248m	3.974	ng/ul	
26) Benzo(g,h,i)perylene	26.21	276	662119	10.052	ng/ul#	93

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

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