

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096863.D
 Acq On : 6 Jul 2018 21:13
 Operator : SJ/JU
 Sample : J3721-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H18

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:09 PM

Quant Time: Jul 09 05:29:04 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	4013	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	21555	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	13692	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	33350m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	31793m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	28254m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	9333	0.26	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	24912m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.19	128	1320	0.025	ng/ul#	89
7) Acenaphthylene	13.74	152	8076	0.144	ng/ul	97
9) Fluorene	15.08	166	1582	0.033	ng/ul	89
12) Phenanthrene	16.82	178	22116m	0.257	ng/ul	
13) Anthracene	16.91	178	13498m	0.185	ng/ul	
15) Fluoranthene	18.85	202	131072m	1.431	ng/ul	
17) Pyrene	19.21	202	182962	1.753	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	119162m	1.589	ng/ul	
19) Chrysene	21.02	228	222406	2.289	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	209956m	2.723	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	71761m	0.852	ng/ul	
23) Benzo(a)pyrene	23.12	252	116665m	1.789	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	65166m	0.962	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	16062m	0.299	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	53302m	0.811	ng/ul	

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

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