

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096861.D
 Acq On : 6 Jul 2018 19:58
 Operator : SJ/JU
 Sample : J3721-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H15

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:28 PM

Quant Time: Jul 09 13:02:26 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	3767	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	19577	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	12617	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	32385m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	31124m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	26012m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	11218	0.35	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	32212m	0.40	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	13.74	152	4636	0.090	ng/ul	99
9) Fluorene	15.08	166	919	0.021	ng/ul	90
12) Phenanthrene	16.82	178	9748m	0.117	ng/ul	
13) Anthracene	16.91	178	8267m	0.117	ng/ul	
15) Fluoranthene	18.85	202	80437m	0.904	ng/ul	
17) Pyrene	19.21	202	93286	0.913	ng/ul#	82
18) Benzo(a)anthracene	20.96	228	66475m	0.906	ng/ul	
19) Chrysene	21.02	228	64139	0.674	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	102060m	1.438	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	37500m	0.484	ng/ul	
23) Benzo(a)pyrene	23.12	252	76658	1.277	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.50	276	29106	0.467	ng/ul#	74
25) Dibenzo(a,h)anthracene	25.51	278	6850	0.138	ng/ul	93
26) Benzo(g,h,i)perylene	26.20	276	24423	0.404	ng/ul#	93

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

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