

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096871.D
 Acq On : 7 Jul 2018 2:11
 Operator : SJ/JU
 Sample : J3765-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F1H08

Manual Integrations
 APPROVED

Sohil
 7/9/2018 4:14:16 PM

Quant Time: Jul 09 05:43:09 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	3436	0.40	ng/ul	0.00
2) Naphthalene-d8	10.14	136	18289	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	12291	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	28818m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	28348m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	27110m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	8184	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	22063m	0.31	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.19	128	10246	0.232	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	5741	0.191	ng/ul	98
7) Acenaphthylene	13.74	152	21049	0.418	ng/ul	97
8) Acenaphthene	14.09	153	1650	0.040	ng/ul#	91
9) Fluorene	15.08	166	13123	0.306	ng/ul	90
12) Phenanthrene	16.82	178	60835m	0.819	ng/ul	
13) Anthracene	16.91	178	413925m	6.560	ng/ul	
15) Fluoranthene	18.85	202	209418m	2.646	ng/ul	
17) Pyrene	19.21	202	279410	3.003	ng/ul#	81
18) Benzo(a)anthracene	20.96	228	167292m	2.502	ng/ul	
19) Chrysene	21.02	228	251625	2.905	ng/ul	99
21) Benzo(b)fluoranthene	22.54	252	557250m	7.531	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	139133m	1.722	ng/ul	
23) Benzo(a)pyrene	23.12	252	252483m	4.034	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	206021m	3.171	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	49355m	0.956	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	168688m	2.675	ng/ul	

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

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