

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE103018\
 Data File : BE098110.D
 Acq On : 30 Oct 2018 12:15
 Operator : SJ/JU
 Sample : J5598-16DL 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE6DL

Manual Integrations
 APPROVED

Sohil
 10/30/2018 5:00:46 PM

Quant Time: Oct 30 13:02:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 30 01:14:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.87	152	1088	0.40	ng/ul	0.00
2) Naphthalene-d8	10.67	136	5464	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3932	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	9567	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	13629	0.40	ng/ul	0.00
20) Perylene-d12	24.02	264	10769	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	185	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	551	0.02	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.72	128	2100	0.161	ng/ul	96
5) 2-Methylnaphthalene	12.34	142	682	0.069	ng/ul	99
7) Acenaphthylene	14.25	152	1821	0.117	ng/ul#	94
12) Phenanthrene	17.33	178	4773	0.216	ng/ul	99
13) Anthracene	17.42	178	3699	0.170	ng/ul	99
15) Fluoranthene	19.35	202	24674	0.876	ng/ul	99
17) Pyrene	19.71	202	31302	0.984	ng/ul	100
18) Benzo(a)anthracene	21.45	228	21463	0.536	ng/ul	98
19) Chrysene	21.50	228	33124	1.069	ng/ul	99
21) Benzo(b)fluoranthene	23.24	252	64317m	2.156	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	18989m	0.650	ng/ul	
23) Benzo(a)pyrene	23.90	252	18572	0.647	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.70	276	23361	0.710	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.73	278	5758	0.210	ng/ul#	90
26) Benzo(g,h,i)perylene	27.54	276	19898	0.719	ng/ul	98

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

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