

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098086.D
 Acq On : 26 Oct 2018 13:03
 Operator : SJ/JU
 Sample : J5598-15
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE3

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:19 PM

Quant Time: Oct 27 00:00:14 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 : Fri Oct 26 23:15:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1122	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5593	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3964	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	9464m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	10594m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	10394m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1983	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	7243m	0.24	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.72	128	1339	0.100	ng/ul#	93
5) 2-Methylnaphthalene	12.34	142	423	0.042	ng/ul	94
7) Acenaphthylene	14.26	152	5278	0.337	ng/ul	99
8) Acenaphthene	14.60	153	358	0.031	ng/ul#	89
9) Fluorene	15.59	166	405	0.029	ng/ul#	92
12) Phenanthrene	17.33	178	12028m	0.551	ng/ul	
13) Anthracene	17.42	178	7784m	0.361	ng/ul	
15) Fluoranthene	19.35	202	123021m	4.418	ng/ul	
17) Pyrene	19.71	202	156598	6.331	ng/ul	99
18) Benzo(a)anthracene	21.45	228	148866m	4.779	ng/ul	
19) Chrysene	21.51	228	147280	6.112	ng/ul	99
21) Benzo(b)fluoranthene	23.25	252	228670m	7.940	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	63550m	2.254	ng/ul	
23) Benzo(a)pyrene	23.91	252	76024	2.745	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.72	276	62098	1.957	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.74	278	17343	0.656	ng/ul	99
26) Benzo(g,h,i)perylene	27.55	276	46313	1.733	ng/ul	97

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

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