

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102618\
 Data File : BE098085.D
 Acq On : 26 Oct 2018 12:27
 Operator : SJ/JU
 Sample : J5598-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AE2

Manual Integrations
 APPROVED

Sohil
 10/29/2018 3:09:18 PM

Quant Time: Oct 26 23:54:40 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	7.87	152	1094	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	5476	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	4003	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	9346m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	10523m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	12220m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	1794	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	6118m	0.20	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.73	128	13632	1.040	ng/ul	98
5) 2-Methylnaphthalene	12.34	142	4813	0.486	ng/ul	100
7) Acenaphthylene	14.26	152	39221	2.483	ng/ul	97
8) Acenaphthene	14.60	153	2675	0.227	ng/ul	95
9) Fluorene	15.59	166	2481	0.175	ng/ul#	98
12) Phenanthrene	17.33	178	55863m	2.590	ng/ul	
13) Anthracene	17.42	178	48445m	2.274	ng/ul	
15) Fluoranthene	19.36	202	910818m	33.119	ng/ul	
17) Pyrene	19.72	202	1289580	52.485	ng/ul	98
18) Benzo(a)anthracene	21.46	228	1061119m	34.296	ng/ul	
19) Chrysene	21.51	228	928089	38.777	ng/ul	98
21) Benzo(b)fluoranthene	23.27	252	1858088m	54.880	ng/ul	
22) Benzo(k)fluoranthene	23.30	252	516140m	15.574	ng/ul	
23) Benzo(a)pyrene	23.93	252	653930m	20.082	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.74	276	514598m	13.791	ng/ul	
25) Dibenzo(a,h)anthracene	26.75	278	135289m	4.350	ng/ul	
26) Benzo(g,h,i)perylene	27.57	276	384384m	12.237	ng/ul	

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

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