

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097466.D
 Acq On : 13 Sep 2018 00:35
 Operator : SJ/JU
 Sample : J4792-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ0

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:33 PM

Quant Time: Sep 13 03:48:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 13 01:51:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1034	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5964	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4310	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	10479m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11345m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	12163m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	1909	0.21	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	6220m	0.18	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.17	128	9644	0.663	ng/ul#	92
5) 2-Methylnaphthalene	11.79	142	3653	0.397	ng/ul	98
7) Acenaphthylene	13.72	152	9457	0.602	ng/ul	99
8) Acenaphthene	14.07	153	6478	0.480	ng/ul	97
9) Fluorene	15.07	166	8011	0.518	ng/ul	93
12) Phenanthrene	16.80	178	202838m	7.520	ng/ul	
13) Anthracene	16.89	178	39514m	1.742	ng/ul	
15) Fluoranthene	18.83	202	424867m	12.260	ng/ul	
17) Pyrene	19.20	202	387877	12.705	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	209384m	7.185	ng/ul	
19) Chrysene	21.00	228	190305	5.668	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	211626m	6.497	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	74068m	2.039	ng/ul	
23) Benzo(a)pyrene	23.11	252	148628m	4.932	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	91731m	2.349	ng/ul	
25) Dibenzo(a,h)anthracene	25.51	278	26774m	0.809	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	80830	2.428	ng/ul#	92

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

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