

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097468.D
 Acq On : 13 Sep 2018 1:49
 Operator : SJ/JU
 Sample : J4792-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ3

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:36 PM

Quant Time: Sep 13 04:11:45 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	1152	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	6565	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4611	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	10985m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11598m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	12163m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	2535	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	8458m	0.23	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.17	128	17776	1.110	ng/ul#	92
5) 2-Methylnaphthalene	11.79	142	6328	0.626	ng/ul	98
7) Acenaphthylene	13.73	152	2529	0.151	ng/ul#	85
8) Acenaphthene	14.07	153	22483	1.556	ng/ul	96
9) Fluorene	15.07	166	29135	1.761	ng/ul	92
12) Phenanthrene	16.80	178	465016m	16.445	ng/ul	
13) Anthracene	16.89	178	114693m	4.824	ng/ul	
15) Fluoranthene	18.84	202	691633m	19.039	ng/ul	
17) Pyrene	19.20	202	598232	19.168	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	288734m	9.691	ng/ul	
19) Chrysene	21.01	228	266892	7.775	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	309521m	9.503	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	109010m	3.001	ng/ul	
23) Benzo(a)pyrene	23.11	252	232016m	7.700	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	148010	3.790	ng/ul#	71
25) Dibenzo(a,h)anthracene	25.50	278	39374m	1.190	ng/ul	
26) Benzo(g,h,i)perylene	26.20	276	127908	3.841	ng/ul#	90

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

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