

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097459.D
 Acq On : 12 Sep 2018 20:15
 Operator : SJ/JU
 Sample : J4792-20
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z35

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:24 PM

Quant Time: Sep 13 02:29:33 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	899	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4989	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3569	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	9178m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10343m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9986m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	1686	0.22	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	6372m	0.21	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.17	128	432	0.036	ng/ul#	75
7) Acenaphthylene	13.72	152	442	0.034	ng/ul#	69
8) Acenaphthene	14.07	153	479	0.043	ng/ul#	93
9) Fluorene	15.06	166	532	0.042	ng/ul#	88
12) Phenanthrene	16.80	178	11527m	0.488	ng/ul	
13) Anthracene	16.89	178	2485m	0.125	ng/ul	
15) Fluoranthene	18.83	202	23170m	0.763	ng/ul	
17) Pyrene	19.20	202	22998	0.826	ng/ul#	80
18) Benzo(a)anthracene	20.95	228	11905m	0.448	ng/ul	
19) Chrysene	21.00	228	11265	0.368	ng/ul	94
21) Benzo(b)fluoranthene	22.53	252	14639m	0.547	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	3922m	0.132	ng/ul	
23) Benzo(a)pyrene	23.10	252	10074m	0.407	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.49	276	6583m	0.205	ng/ul	
25) Dibenzo(a,h)anthracene	25.50	278	1654m	0.061	ng/ul	
26) Benzo(g,h,i)perylene	26.19	276	5496m	0.201	ng/ul	

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

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