

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097460.D
 Acq On : 12 Sep 2018 20:51
 Operator : SJ/JU
 Sample : J4792-22
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z38

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:26 PM

Quant Time: Sep 13 02:35:21 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	888	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5004	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3686	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	9044m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10369m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10106m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2600	0.33	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	9180m	0.31	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	16.80	178	1428m	0.061	ng/ul	
15) Fluoranthene	18.83	202	3371m	0.113	ng/ul	
17) Pyrene	19.20	202	3138m	0.112	ng/ul	
18) Benzo(a)anthracene	20.95	228	1882m	0.071	ng/ul	
19) Chrysene	21.01	228	2135m	0.070	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	2982m	0.110	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	848m	0.028	ng/ul	
23) Benzo(a)pyrene	23.10	252	1717	0.069	ng/ul#	15
24) Indeno(1,2,3-cd)pyrene	25.50	276	1339	0.041	ng/ul#	87
26) Benzo(g,h,i)perylene	26.19	276	1367	0.049	ng/ul#	68

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

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