

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097475.D
 Acq On : 13 Sep 2018 6:09
 Operator : SJ/JU
 Sample : J4792-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YY4

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:44 PM

Quant Time: Sep 13 08:20:31 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 07:48:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	1137	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	6617	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4931	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	12334m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12575m	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	13469m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2759	0.27	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9758m	0.24	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.17	128	5559	0.345	ng/ul#	93
5) 2-Methylnaphthalene	11.79	142	2950	0.289	ng/ul	97
7) Acenaphthylene	13.73	152	6300	0.351	ng/ul	97
8) Acenaphthene	14.07	153	4291	0.278	ng/ul	96
9) Fluorene	15.07	166	4398	0.249	ng/ul	89
12) Phenanthrene	16.80	178	68268m	2.150	ng/ul	
13) Anthracene	16.90	178	14158m	0.530	ng/ul	
15) Fluoranthene	18.84	202	144130m	3.534	ng/ul	
17) Pyrene	19.20	202	148904m	4.400	ng/ul	
18) Benzo(a)anthracene	20.96	228	77088m	2.386	ng/ul	
19) Chrysene	21.01	228	96457m	2.592	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	130307m	3.613	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	34300m	0.853	ng/ul	
23) Benzo(a)pyrene	23.12	252	78564	2.354	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.52	276	68210	1.577	ng/ul#	73
25) Dibenzo(a,h)anthracene	25.53	278	18722	0.511	ng/ul#	65
26) Benzo(g,h,i)perylene	26.22	276	64489	1.749	ng/ul	94

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

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