

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN120720\
 Data File : BN013037.D
 Acq On : 10 Dec 2020 09:03
 Operator : CG/JU
 Sample : L4941-05DL 2X
 Misc :
 ALS Vial : 107 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 C0AY4DL

Quant Time: Dec 10 10:28:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 07:33:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	1299	0.40	ng/ul	-0.03
2) Naphthalene-d8	10.13	136	5149	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.04	164	2846	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.80	188	5708	0.40	ng/ul	0.00
16) Chrysene-d12	21.03	240	4043	0.40	ng/ul	0.00
20) Perylene-d12	23.13	264	3742	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	521	0.07	ng/ul	-0.02
14) Fluoranthene-d10	18.84	212	1154	0.07	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	13.75	152	415	0.031	ng/ul#	77
9) Fluorene	15.10	166	271	0.021	ng/ul#	87
12) Phenanthrene	16.84	178	7798	0.407	ng/ul	99
13) Anthracene	16.93	178	969	0.059	ng/ul#	86
15) Fluoranthene	18.87	202	20991	0.923	ng/ul	100
17) Pyrene	19.24	202	18088	0.947	ng/ul	99
18) Benzo(a)anthracene	21.01	228	7428	0.444	ng/ul	99
19) Chrysene	21.07	228	10255	0.555	ng/ul	99
21) Benzo(b)fluoranthene	22.51	252	18189	1.051	ng/ul	93
22) Benzo(k)fluoranthene	22.51	252	18189	0.951	ng/ul	94
23) Benzo(a)pyrene	22.95	252	3976	0.248	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.18	276	5627	0.267	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.19	278	1320	0.079	ng/ul#	67
26) Benzo(g,h,i)perylene	25.81	276	4880	0.256	ng/ul	97

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

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