

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102120\
 Data File : BN012382.D
 Acq On : 21 Oct 2020 13:54
 Operator : CG/JU
 Sample : L4425-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2P78

Quant Time: Oct 21 14:23:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 21 12:12:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	726	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.36	136	3059	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	1813	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	3549	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	3072	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	3165	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	1517	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	3364	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	2244	0.254	ng/ul#	91
5) 2-Methylnaphthalene	12.03	142	3164	0.564	ng/ul	99
7) Acenaphthylene	13.95	152	1104	0.140	ng/ul#	86
8) Acenaphthene	14.29	153	2384	0.359	ng/ul	98
9) Fluorene	15.29	166	3832	0.518	ng/ul	99
11) Pentachlorophenol	16.64	266	1182	1.153	ng/ul	100
12) Phenanthrene	17.03	178	4393	0.388	ng/ul	98
13) Anthracene	17.12	178	3599	0.371	ng/ul	98
15) Fluoranthene	19.05	202	5143	0.389	ng/ul	99
17) Pyrene	19.41	202	5111	0.369	ng/ul	99
18) Benzo(a)anthracene	21.17	228	5235	0.460	ng/ul	98
19) Chrysene	21.22	228	6339	0.457	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	8218	0.643	ng/ul	97
22) Benzo(k)fluoranthene	22.76	252	4494	0.318	ng/ul#	82
23) Benzo(a)pyrene	23.27	252	3389	0.273	ng/ul#	69
25) Dibenzo(a,h)anthracene	25.54	278	9253	0.724	ng/ul	97
26) Benzo(a,h,i)perylene	26.19	276	4075	0.278	ng/ul	96

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

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