

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003133.D
 Acq On : 12 Oct 2018 08:56
 Operator : MJ/SJ
 Sample : J5234-17DL 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AKODL

Quant Time: Oct 12 13:01:08 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1581	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.39	136	6826	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	4093	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8573	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	7631	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5956	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.99	152	627	0.06	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	2381	0.10	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.44	128	5667	0.324	ng/ul	98
5) 2-Methylnaphthalene	12.06	142	10479	0.869	ng/ul	96
7) Acenaphthylene	13.98	152	1016	0.054	ng/ul#	34
8) Acenaphthene	14.32	153	1810	0.127	ng/ul	95
9) Fluorene	15.31	166	8856	0.544	ng/ul#	97
12) Phenanthrene	17.06	178	18828	0.788	ng/ul	97
13) Anthracene	17.15	178	5715	0.253	ng/ul#	91
15) Fluoranthene	19.08	202	7037	0.251	ng/ul	92
17) Pyrene	19.45	202	11260	0.380	ng/ul	99
18) Benzo(a)anthracene	21.21	228	3074	0.125	ng/ul#	66
19) Chrysene	21.26	228	2970	0.119	ng/ul#	68
21) Benzo(b)fluoranthene	22.78	252	1861	0.087	ng/ul#	25
22) Benzo(k)fluoranthene	22.78	252	1861	0.084	ng/ul#	26
23) Benzo(a)pyrene	23.34	252	1360	0.071	ng/ul#	34
24) Indeno(1,2,3-cd)pyrene	25.66	276	642	0.029	ng/ul#	49
26) Benzo(g,h,i)perylene	26.34	276	518	0.028	ng/ul#	27

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

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