

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003130.D
 Acq On : 12 Oct 2018 07:10
 Operator : MJ/SJ
 Sample : J5234-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK9

Quant Time: Oct 12 12:25:32 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	7.62	152	1915	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7626	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4856	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	315396	0.40	ng/ul	0.08
16) Chrysene-d12	21.33	240	176	0.40	ng/ul	0.10
20) Perylene-d12	23.55	264	196	0.40	ng/ul	0.10

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.98	152	2873	0.24	ng/ul	-0.03
14) Fluoranthene-d10	19.09	212	4379	0.00	ng/ul	0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.44	128	9959	0.509	ng/ul#	96
5) 2-Methylnaphthalene	12.06	142	36111	2.681	ng/ul	97
7) Acenaphthylene	13.98	152	4873	0.219	ng/ul#	46
8) Acenaphthene	14.32	153	9015	0.535	ng/ul	98
9) Fluorene	15.31	166	40989	2.123	ng/ul	98
12) Phenanthrene	17.05	178	70252	0.080	ng/ul	97
17) Pyrene	19.45	202	54504	79.734	ng/ul	98
18) Benzo(a)anthracene	21.26	228	17935	31.667	ng/ul#	78
19) Chrysene	21.38	228	2165	3.765	ng/ul#	1
21) Benzo(b)fluoranthene	22.78	252	24014	34.020	ng/ul	90
22) Benzo(k)fluoranthene	22.95	252	2947	4.027	ng/ul#	38
23) Benzo(a)pyrene	23.48	252	3754	5.934	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	25.65	276	8958	12.323	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.75	278	710	1.220	ng/ul#	1
26) Benzo(g,h,i)perylene	26.32	276	8164	13.318	ng/ul	92

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

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