

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003131.D
 Acq On : 12 Oct 2018 07:45
 Operator : MJ/SJ
 Sample : J5234-11DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK9DL

Quant Time: Oct 12 12:40:19 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	1576	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.39	136	5889	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	2905	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	4974	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	4648	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5024	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.99	152	458	0.05	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	1514	0.11	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.44	128	1818	0.120	ng/ul#	87
5) 2-Methylnaphthalene	12.07	142	5760	0.554	ng/ul	98
7) Acenaphthylene	13.98	152	928	0.070	ng/ul#	43
8) Acenaphthene	14.32	153	1445	0.143	ng/ul	96
9) Fluorene	15.32	166	6470	0.560	ng/ul#	98
12) Phenanthrene	17.06	178	11717	0.845	ng/ul#	95
13) Anthracene	17.15	178	3593	0.274	ng/ul#	87
15) Fluoranthene	19.08	202	8665	0.533	ng/ul	94
17) Pyrene	19.45	202	10723	0.594	ng/ul	98
18) Benzo(a)anthracene	21.21	228	3884	0.260	ng/ul#	74
19) Chrysene	21.26	228	3854	0.254	ng/ul#	77
21) Benzo(b)fluoranthene	22.78	252	5250	0.290	ng/ul	68
22) Benzo(k)fluoranthene	22.78	252	5250	0.280	ng/ul#	69
23) Benzo(a)pyrene	23.34	252	3083	0.190	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.65	276	1948	0.105	ng/ul	91
25) Dibenzo(a,h)anthracene	25.65	278	468	0.031	ng/ul#	1
26) Benzo(g,h,i)perylene	26.33	276	1819	0.116	ng/ul#	74

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

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