

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101018\
 Data File : BN003060.D
 Acq On : 10 Oct 2018 10:45
 Operator : MJ/SJ
 Sample : J5115-14DL 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC66DL

Quant Time: Oct 10 12:28:17 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 : Wed Oct 10 08:36:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1415	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	6311	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3856	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9165	0.40	ng/ul	-0.01
16) Chrysene-d12	21.22	240	9222	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	8007	0.40	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.03	152	259	0.03	ng/ul	0.02
14) Fluoranthene-d10	19.05	212	734	0.03	ng/ul	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.46	128	449	0.028	ng/ul#	60
8) Acenaphthene	14.33	153	826	0.062	ng/ul	98
9) Fluorene	15.33	166	772	0.050	ng/ul#	95
12) Phenanthrene	17.06	178	14229	0.557	ng/ul	100
13) Anthracene	17.15	178	2952	0.122	ng/ul#	93
15) Fluoranthene	19.09	202	30972	1.033	ng/ul	99
17) Pyrene	19.45	202	28053	0.783	ng/ul	98
18) Benzo(a)anthracene	21.21	228	13642	0.460	ng/ul	98
19) Chrysene	21.26	228	27764	0.921	ng/ul	100
21) Benzo(b)fluoranthene	22.78	252	39492	1.370	ng/ul	96
22) Benzo(k)fluoranthene	22.78	252	39494	1.321	ng/ul	97
23) Benzo(a)pyrene	23.34	252	11625	0.450	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.65	276	10509	0.354	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.65	278	2677	0.113	ng/ul#	73
26) Benzo(g,h,i)perylene	26.33	276	11495	0.459	ng/ul	97

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

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