

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN100318\
 Data File : BN002896.D
 Acq On : 03 Oct 2018 12:22
 Operator : MJ/SJ
 Sample : J5123-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0B46

Quant Time: Oct 03 13:19:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 29 06:35:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	140884	20.00	ng/ul	0.00
18) Naphthalene-d8	10.39	136	601915	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.26	164	367571	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.02	188	800032	20.00	ng/ul	0.00
77) Chrysene-d12	21.22	240	868407	20.00	ng/ul	0.00
85) Perylene-d12	23.42	264	866466	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	0.00	67	0	0.00	ng/ul	
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
46) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	0.00	176	0d	0.00	ng/ul	
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	0.00	188	0d	0.00	ng/ul	
78) Pyrene-d10	0.00	212	0d	0.00	ng/ul	
89) Benzo(a)pyrene-d12	0.00	264	0d	0.00	ng/ul	

Target Compounds

				Ovalue		
69) Phenanthrene	17.06	178	45883	1.041	ng/ul#	91
76) Fluoranthene	19.09	202	141581	3.011	ng/ul	99
79) Pyrene	19.45	202	63428	1.105	ng/ul#	94
82) Benzo(a)anthracene	21.20	228	55547	1.028	ng/ul	98
84) Chrysene	21.26	228	80999	1.600	ng/ul	97
87) Benzo(b)fluoranthene	22.77	252	97820	1.849	ng/ul#	96

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

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