

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121418\
 Data File : BN003937.D
 Acq On : 14 Dec 2018 18:31
 Operator : JU/SJ
 Sample : J6270-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E00

Quant Time: Dec 14 23:11:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN121218MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 14 22:50:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	36765	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	175745	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.47	164	106500	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.21	188	254684	20.00	ng/ul	0.00
77) Chrysene-d12	21.40	240	313482	20.00	ng/ul	0.00
85) Perylene-d12	23.72	264	291616	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.29	96	2168	3.86	ng/uL	0.00
5) Phenol-d5	6.99	99	76685	21.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.16	67	41558	21.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.36	132	63564	22.44	ng/ul	0.00
13) 4-Methylphenol-d8	8.53	113	62750	20.72	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	32584	23.40	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	37317	23.66	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.25	165	67457	23.32	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	51452	19.01	ng/ul	0.00
43) Dimethylphthalate-d6	13.87	166	208744	21.96	ng/ul	0.00
46) Acenaphthylene-d8	14.16	160	274337	24.32	ng/ul	0.00
51) 4-Nitrophenol-d4	14.67	143	35651	18.14	ng/ul	0.00
57) Fluorene-d10	15.46	176	195290	23.76	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.58	200	15211	8.29	ng/ul	0.00
70) Anthracene-d10	17.31	188	318227	24.73	ng/ul	0.00
78) Pyrene-d10	19.60	212	378145	26.96	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.56	264	396639	25.07	ng/ul	0.00

Target Compounds				Ovalue		
6) Phenol	7.02	94	18014	4.938	ng/ul	98
44) Dimethylphthalate	13.92	163	68862	7.420	ng/ul	99
76) Fluoranthene	19.26	202	21332	1.183	ng/ul	98
79) Pyrene	19.63	202	24747	1.413	ng/ul	97
84) Chrysene	21.43	228	21120	1.153	ng/ul	97
87) Benzo(b)fluoranthene	23.01	252	39240	2.246	ng/ul#	95
93) Benzo(a,h,i)perylene	26.79	276	21645	1.179	ng/ul	95

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

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