

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM100319\
 Data File : BM022956.D
 Acq On : 04 Oct 2019 14:16
 Operator : JU
 Sample : K5006-08
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETJW4

Quant Time: Oct 04 16:54:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 27 18:03:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	186856	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.58	136	810087	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.43	164	475164	20.00	ng/ul	-0.02
62) Phenanthrene-d10	17.17	188	822008	20.00	ng/ul	-0.02
78) Chrysene-d12	21.34	240	681358	20.00	ng/ul	-0.02
86) Perylene-d12	23.61	264	784325	20.00	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.26	96	21971	5.09	ng/uL	0.00
5) Phenol-d5	6.96	99	127158	7.70	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	7.12	67	264557	29.28	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.33	132	331439	24.93	ng/ul	-0.02
13) 4-Methylphenol-d8	8.50	113	219688	16.25	ng/ul	-0.02
19) Nitrobenzene-d5	8.95	128	189203	30.26	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.67	143	205713	30.13	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.20	165	367423	27.03	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.72	131	42170	2.55	ng/ul	-0.02
44) Dimethylphthalate-d6	13.84	166	1117676	30.31	ng/ul	-0.02
47) Acenaphthylene-d8	14.12	160	1285103	30.18	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.63	143	35421	4.86	ng/ul	-0.01
58) Fluorene-d10	15.42	176	925140	29.43	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.53	200	126830	23.71	ng/ul	-0.02
71) Anthracene-d10	17.26	188	1401121	34.90	ng/ul	-0.02
79) Pyrene-d10	19.56	212	1361920	37.14	ng/ul	-0.02
90) Benzo(a)pyrene-d12	23.47	264	1443903	35.59	ng/ul	-0.03

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

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