

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102319\
 Data File : BM023377.D
 Acq On : 24 Oct 2019 13:45
 Operator : JU
 Sample : AR1268
 Misc : GCMS CONFIRMATION
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 24 15:03:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 24 13:46:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	536209	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.41	136	2060476	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.28	164	1180090	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.02	188	2235390	20.00	ng/ul	-0.01
78) Chrysene-d12	21.22	240	1886710	20.00	ng/ul	-0.01
86) Perylene-d12	23.42	264	2348997	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	160	0.01	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.99	67	147	0.01	ng/ul	0.00
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	10.42	131	2800	0.07	ng/ul	-0.14
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	13.97	160	118	0.00	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.27	176	1968	0.03	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	17.02	188	2235390	21.16	ng/ul	-0.11
79) Pyrene-d10	19.42	212	933	0.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.42	264	2326812	19.44	ng/ul	0.13

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

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