

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102919\
 Data File : BM023443.D
 Acq On : 29 Oct 2019 15:21
 Operator : JU
 Sample : K5423-10
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Oct 29 18:05:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 25 07:44:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	483552	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.40	136	1941411	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.27	164	1144551	20.00	ng/ul	-0.01
62) Phenanthrene-d10	17.02	188	2259624	20.00	ng/ul	-0.01
78) Chrysene-d12	21.22	240	2268959	20.00	ng/ul	-0.01
86) Perylene-d12	23.41	264	2664600	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.15	96	65394	6.42	ng/uL	0.00
5) Phenol-d5	6.80	99	324869	8.00	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.96	67	769916	32.34	ng/ul	-0.02
9) 2-Chlorophenol-d4	7.16	132	885490	27.89	ng/ul	-0.01
13) 4-Methylphenol-d8	8.33	113	568464	17.23	ng/ul	-0.02
19) Nitrobenzene-d5	8.77	128	527012	36.06	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.49	143	581686	35.87	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	10.03	165	972997	30.53	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.52	131	17957	0.49	ng/ul	-0.04
44) Dimethylphthalate-d6	13.69	166	3077783	35.41	ng/ul	-0.01
47) Acenaphthylene-d8	13.96	160	3545733	35.61	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.49	143	102956	6.84	ng/ul	0.00
58) Fluorene-d10	15.27	176	2586581	35.14	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.40	200	401412	28.53	ng/ul	-0.01
71) Anthracene-d10	17.12	188	4038196	37.82	ng/ul	-0.02
79) Pyrene-d10	19.42	212	4504027	36.10	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.27	264	5278162	38.88	ng/ul	-0.01

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

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