

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101519\
 Data File : BM023183.D
 Acq On : 16 Oct 2019 01:09
 Operator : JU
 Sample : PB123737BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK37

Quant Time: Oct 16 06:50:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 02:28:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	456267	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	1985909	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	1288816	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	2694908	20.00	ng/ul	-0.01
78) Chrysene-d12	21.26	240	2341088	20.00	ng/ul	-0.01
86) Perylene-d12	23.47	264	2309482	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.20	96	66941	6.46	ng/uL	0.00
5) Phenol-d5	6.86	99	1170961	29.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.02	67	733757	32.64	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	924581	30.40	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	952092	29.83	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	441947	32.89	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	440043	34.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	956405	29.40	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	1386605	35.03	ng/ul	0.00
44) Dimethylphthalate-d6	13.74	166	3222469	32.63	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	3714517	32.07	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.52	143	449298	28.10	ng/ul	0.00
58) Fluorene-d10	15.32	176	2694367	32.02	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.43	200	264947	24.80	ng/ul	-0.01
71) Anthracene-d10	17.17	188	4288981	33.01	ng/ul	-0.01
79) Pyrene-d10	19.46	212	4527898	39.94	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.33	264	3930982	33.58	ng/ul	-0.02

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

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