

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000733.D
 Acq On : 15 Oct 2019 00:56
 Operator : JU
 Sample : PB123959BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK59

Quant Time: Oct 15 04:53:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 03:51:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	180136	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	700347	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	391159	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	724752	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	616645	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	642183	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.36	96	36318	8.36	ng/uL	0.00
5) Phenol-d5	6.39	99	576110	41.46	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	287407	41.37	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	486295	41.54	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	450303	42.42	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	228420	42.91	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	247918	43.73	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	427204	39.37	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	611821	50.70	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1188990	43.42	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1406334	41.75	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	170085	40.33	ng/ul	0.00
58) Fluorene-d10	10.32	176	979972	41.83	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	109971	31.86	ng/ul	0.00
71) Anthracene-d10	11.35	188	1437635	42.82	ng/ul	0.00
79) Pyrene-d10	12.73	212	1482897	42.53	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1368756	43.17	ng/ul	0.00

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

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