

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000751.D
 Acq On : 15 Oct 2019 08:33
 Operator : JU
 Sample : PB123736BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 15 09:14:20 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	244386	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	951069	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	561430	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	1141916	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	1051595	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	1085842	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.38	96	25	0.00	ng/uL	0.01
5) Phenol-d5	6.38	99	9	0.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	23	0.00	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	18	0.00	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	34	0.00	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	42	0.01	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	14	0.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.83	165	26	0.00	ng/ul	-0.06
29) 4-Chloroaniline-d4	8.08	131	11	0.00	ng/ul	-0.01
44) Dimethylphthalate-d6	9.50	166	15	0.00	ng/ul	0.00
47) Acenaphthylene-d8	9.80	160	255031	5.28	ng/ul	0.15
52) 4-Nitrophenol-d4	9.92	143	7	0.00	ng/ul	0.00
58) Fluorene-d10	10.32	176	1031	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.40	200	11	0.00	ng/ul	0.00
71) Anthracene-d10	11.35	188	6122	0.12	ng/ul	0.00
79) Pyrene-d10	12.75	212	27	0.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.45	264	1084006	20.22	ng/ul	0.08

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

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