

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000749.D
 Acq On : 15 Oct 2019 07:45
 Operator : JU
 Sample : PB123845BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 15 08:13:32 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	266784	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	1044266	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	615215	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	1221989	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	1058426	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	1042809	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	96	0.01	ng/uL	0.01
5) Phenol-d5	6.38	99	11	0.00	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.42	67	50	0.00	ng/ul	-0.01
9) 2-Chlorophenol-d4	6.52	132	28	0.00	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	13	0.00	ng/ul	0.00
19) Nitrobenzene-d5	7.32	128	22	0.00	ng/ul	0.00
22) 2-Nitrophenol-d4	7.64	143	22	0.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.81	165	52	0.00	ng/ul	-0.08
29) 4-Chloroaniline-d4	8.10	131	33	0.00	ng/ul	0.01
44) Dimethylphthalate-d6	9.49	166	29	0.00	ng/ul	0.00
47) Acenaphthylene-d8	9.80	160	277720	5.24	ng/ul	0.15
52) 4-Nitrophenol-d4	9.91	143	13	0.00	ng/ul	0.00
58) Fluorene-d10	10.32	176	1203	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.40	200	8	0.00	ng/ul	0.00
71) Anthracene-d10	11.35	188	7737	0.14	ng/ul	0.00
79) Pyrene-d10	12.74	212	43	0.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.45	264	1041544	20.23	ng/ul	0.08

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

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