

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101519\
 Data File : BP000748.D
 Acq On : 15 Oct 2019 07:21
 Operator : JU
 Sample : PB123889BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 15 08:01:08 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	181399	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	696113	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	420016	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	871024	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	813016	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	809611	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.37	96	25	0.01	ng/uL	0.00
5) Phenol-d5	6.39	99	6	0.00	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	19	0.00	ng/ul	-0.01
9) 2-Chlorophenol-d4	6.53	132	21	0.00	ng/ul	0.00
13) 4-Methylphenol-d8	7.12	113	7	0.00	ng/ul	-0.01
19) Nitrobenzene-d5	7.28	128	13	0.00	ng/ul	-0.03
22) 2-Nitrophenol-d4	7.63	143	8	0.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	17	0.00	ng/ul	0.00
29) 4-Chloroaniline-d4	8.08	131	37	0.00	ng/ul	-0.01
44) Dimethylphthalate-d6	9.49	166	35	0.00	ng/ul	0.00
47) Acenaphthylene-d8	9.80	160	189458	5.24	ng/ul	0.15
52) 4-Nitrophenol-d4	9.91	143	14	0.00	ng/ul	0.00
58) Fluorene-d10	10.32	176	814	0.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	9	0.00	ng/ul	0.00
71) Anthracene-d10	11.35	188	5691	0.14	ng/ul	0.00
79) Pyrene-d10	12.74	212	67	0.00	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.45	264	809611	20.26	ng/ul	0.08

Target Compounds	Ovalue

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

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