

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000926.D
 Acq On : 19 Oct 2019 19:51
 Operator : JU
 Sample : PB123959BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK59

Quant Time: Oct 19 23:06:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 19 22:14:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	123515	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	464654	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	266636	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	518682	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	505826	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	589105	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.33	96	18817	6.32	ng/uL	0.00
5) Phenol-d5	6.38	99	290323	30.47	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	143310	30.09	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	256046	31.90	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	233758	32.11	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	116973	33.12	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	130223	34.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.87	165	229734	31.91	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	315832	39.45	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	635406	34.04	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	744370	32.42	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	84366	29.35	ng/ul	-0.01
58) Fluorene-d10	10.31	176	527456	33.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	56622	22.92	ng/ul	0.00
71) Anthracene-d10	11.34	188	802731	33.41	ng/ul	0.00
79) Pyrene-d10	12.72	212	875387	30.60	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	971340	33.40	ng/ul	0.00

Target Compounds Ovalue

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

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