

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000927.D
 Acq On : 19 Oct 2019 20:15
 Operator : JU
 Sample : PB124128BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK28

Quant Time: Oct 19 23:08:02 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	136788	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	509900	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	285345	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	529962	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	500090	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	589607	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.35	96	4552	1.38	ng/uL	0.01
5) Phenol-d5	6.38	99	74812	7.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	149355	28.31	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	226126	25.44	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	141248	17.52	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	126338	32.59	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	140306	33.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.87	165	233920	29.61	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	386	0.04	ng/ul	0.01
44) Dimethylphthalate-d6	9.48	166	670374	33.56	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	797322	32.45	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	19216	6.25	ng/ul	0.00
58) Fluorene-d10	10.31	176	569827	33.34	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	62541	24.78	ng/ul	0.00
71) Anthracene-d10	11.34	188	865639	35.26	ng/ul	0.00
79) Pyrene-d10	12.72	212	927798	32.81	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1017854	34.97	ng/ul	0.00

Target Compounds Ovalue

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

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