

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP091019\
 Data File : BP000168.D
 Acq On : 10 Sep 2019 16:34
 Operator : HP/JU
 Sample : PB122972BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Sep 11 01:57:08 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP090519MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 02:43:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	402195	20.00	ng/ul	0.00
18) Naphthalene-d8	8.17	136	1592804	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.95	164	883942	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.45	188	1642058	20.00	ng/ul	0.00
78) Chrysene-d12	14.14	240	1286173	20.00	ng/ul	0.00
86) Perylene-d12	15.68	264	1261614	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.61	96	63770	5.62	ng/uL	0.00
5) Phenol-d5	6.51	99	1009995	29.48	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.57	67	522122	29.15	ng/ul	0.00
9) 2-Chlorophenol-d4	6.66	132	834649	30.13	ng/ul	0.00
13) 4-Methylphenol-d8	7.26	113	773586	30.06	ng/ul	0.00
19) Nitrobenzene-d5	7.45	128	385744	31.90	ng/ul	0.00
22) 2-Nitrophenol-d4	7.77	143	390519	33.17	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	8.02	165	709849	28.40	ng/ul	0.00
29) 4-Chloroaniline-d4	8.23	131	1034863	33.49	ng/ul	0.00
44) Dimethylphthalate-d6	9.63	166	1983481	30.39	ng/ul	0.00
47) Acenaphthylene-d8	9.79	160	2368033	29.36	ng/ul	0.00
52) 4-Nitrophenol-d4	10.03	143	329101	28.22	ng/ul	0.00
58) Fluorene-d10	10.47	176	1657727	29.91	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.53	200	175745	23.29	ng/ul	0.00
71) Anthracene-d10	11.50	188	2380448	29.94	ng/ul	0.00
79) Pyrene-d10	12.90	212	2403008	31.41	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.59	264	2070867	31.67	ng/ul	0.00

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

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