

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000580.D
 Acq On : 09 Oct 2019 19:54
 Operator : JU
 Sample : K5201-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AC0

Quant Time: Oct 10 01:21:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 09 17:20:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	228689	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	882948	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	520095	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	984359	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	851727	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	945616	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.36	96	28357	5.14	ng/uL	-0.01
5) Phenol-d5	6.39	99	110344	6.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	231525	26.25	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	320195	21.54	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	184566	13.69	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	184757	27.53	ng/ul	0.00
22) 2-Nitrophenol-d4	7.64	143	201245	28.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.89	165	346872	25.36	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	20243	1.33	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1122674	30.84	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1253004	27.98	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	35160	6.27	ng/ul	0.00
58) Fluorene-d10	10.32	176	921223	29.57	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	119947	25.59	ng/ul	0.00
71) Anthracene-d10	11.35	188	1401534	30.74	ng/ul	0.00
79) Pyrene-d10	12.73	212	1587636	32.96	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1577653	33.79	ng/ul	0.00

Target Compounds					Ovalue	
6) Phenol	6.40	94	24938	1.402	ng/ul	95
45) Dimethylphthalate	9.51	163	83384	2.299	ng/ul	100

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

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