

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000922.D
 Acq On : 19 Oct 2019 18:15
 Operator : JU
 Sample : K5395-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0051

Quant Time: Oct 19 20:29:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 09:27:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	126726	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	468311	20.00	ng/ul	-0.01
36) Acenaphthene-d10	9.79	164	263455	20.00	ng/ul	-0.01
62) Phenanthrene-d10	11.28	188	495228	20.00	ng/ul	-0.02
78) Chrysene-d12	13.96	240	470378	20.00	ng/ul	-0.02
86) Perylene-d12	15.43	264	550950	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.43	96	161	0.05	ng/uL	0.04
5) Phenol-d5	6.38	99	73326	7.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	148291	30.34	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	225765	27.41	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	140396	18.80	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	124339	34.93	ng/ul	-0.01
22) 2-Nitrophenol-d4	7.62	143	138911	36.64	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	7.88	165	232244	32.01	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	17570	2.18	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	655593	35.55	ng/ul	-0.01
47) Acenaphthylene-d8	9.63	160	797354	35.15	ng/ul	-0.01
52) 4-Nitrophenol-d4	9.92	143	18440	6.49	ng/ul	0.00
58) Fluorene-d10	10.31	176	555451	35.20	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	10.39	200	61396	26.03	ng/ul	-0.01
71) Anthracene-d10	11.34	188	840062	36.62	ng/ul	-0.01
79) Pyrene-d10	12.72	212	895789	33.68	ng/ul	-0.02
90) Benzo(a)pyrene-d12	15.35	264	991757	36.46	ng/ul	-0.02

Target Compounds	Ovalue

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

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