

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM011720\
 Data File : BM024505.D
 Acq On : 17 Jan 2020 20:58
 Operator : JU
 Sample : PB126070BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK70

Quant Time: Jan 18 00:30:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jan 18 00:06:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	179864	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	792596	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	607476	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.86	188	1427209	20.00	ng/ul	0.00
78) Chrysene-d12	21.06	240	1460422	20.00	ng/ul	0.00
86) Perylene-d12	23.18	264	1551366	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	26799	7.28	ng/uL	0.00
5) Phenol-d5	6.65	99	429408	32.21	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.81	67	258582	34.92	ng/ul	0.00
9) 2-Chlorophenol-d4	7.00	132	408232	35.85	ng/ul	0.00
13) 4-Methylphenol-d8	8.17	113	401274	32.64	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	216995	38.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.31	143	249183	41.32	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	464554	34.22	ng/ul	0.00
29) 4-Chloroaniline-d4	10.36	131	650493	43.10	ng/ul	0.00
44) Dimethylphthalate-d6	13.53	166	1811408	38.06	ng/ul	0.00
47) Acenaphthylene-d8	13.79	160	2030915	37.07	ng/ul	0.00
52) 4-Nitrophenol-d4	14.32	143	257872	32.46	ng/ul	0.00
58) Fluorene-d10	15.10	176	1553991	37.09	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.23	200	262864	32.26	ng/ul	0.00
71) Anthracene-d10	16.95	188	2484191	37.43	ng/ul	0.00
79) Pyrene-d10	19.26	212	2906355	37.23	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.05	264	2952329	36.69	ng/ul	0.00

Target Compounds	Ovalue

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

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