

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021720\
 Data File : BM024923.D
 Acq On : 18 Feb 2020 09:44
 Operator : CG/JU
 Sample : PB126689BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK89

Quant Time: Feb 18 10:25:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM020620MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 18 02:53:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	388401	20.00	ng/ul	0.00
18) Naphthalene-d8	10.15	136	1397397	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.04	164	821123	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.79	188	1701251	20.00	ng/ul	0.00
78) Chrysene-d12	21.01	240	1695901	20.00	ng/ul	0.00
86) Perylene-d12	23.11	264	1762819	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.04	96	69652	8.63	ng/uL	0.00
5) Phenol-d5	6.59	99	1017594	39.44	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.75	67	619672	42.52	ng/ul	0.00
9) 2-Chlorophenol-d4	6.93	132	865138	37.42	ng/ul	0.00
13) 4-Methylphenol-d8	8.10	113	774534	35.57	ng/ul	0.00
19) Nitrobenzene-d5	8.53	128	391239	38.72	ng/ul	0.00
22) 2-Nitrophenol-d4	9.25	143	452687	37.91	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.78	165	802366	33.31	ng/ul	0.00
29) 4-Chloroaniline-d4	10.29	131	991717	44.02	ng/ul	0.00
44) Dimethylphthalate-d6	13.47	166	2123224	34.26	ng/ul	0.00
47) Acenaphthylene-d8	13.72	160	2682575	37.06	ng/ul	0.00
52) 4-Nitrophenol-d4	14.26	143	316856	31.46	ng/ul	0.00
58) Fluorene-d10	15.04	176	1884164	34.33	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.18	200	315571	28.46	ng/ul	0.00
71) Anthracene-d10	16.89	188	2780760	35.81	ng/ul	0.00
79) Pyrene-d10	19.20	212	3087192	36.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	22.98	264	3246530	36.67	ng/ul	0.00

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

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