

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025803.D
 Acq On : 26 Mar 2020 17:03
 Operator : CG/JU
 Sample : PB127666BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK66

Quant Time: Mar 26 17:37:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 24 00:45:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	229949	20.00	ng/ul	0.00
18) Naphthalene-d8	10.33	136	925109	20.00	ng/ul	-0.01
36) Acenaphthene-d10	14.20	164	568579	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.95	188	1228084	20.00	ng/ul	-0.01
78) Chrysene-d12	21.14	240	1334367	20.00	ng/ul	-0.01
86) Perylene-d12	23.30	264	1492308	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.13	96	39713	7.25	ng/uL	0.00
5) Phenol-d5	6.75	99	630902	34.76	ng/ul	-0.01
7) Bis-(2-Chloroethyl)ether-d	6.91	67	375120	36.21	ng/ul	-0.01
9) 2-Chlorophenol-d4	7.10	132	534432	36.28	ng/ul	-0.01
13) 4-Methylphenol-d8	8.27	113	522030	35.01	ng/ul	-0.01
19) Nitrobenzene-d5	8.71	128	267514	38.91	ng/ul	0.00
22) 2-Nitrophenol-d4	9.43	143	287827	39.37	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.96	165	528358	35.35	ng/ul	0.00
29) 4-Chloroaniline-d4	10.47	131	750016	43.21	ng/ul	-0.01
44) Dimethylphthalate-d6	13.63	166	1634397	37.82	ng/ul	0.00
47) Acenaphthylene-d8	13.89	160	1994981	38.53	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.42	143	281627	34.23	ng/ul	0.00
58) Fluorene-d10	15.20	176	1446873	38.14	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.33	200	224840	30.01	ng/ul	-0.01
71) Anthracene-d10	17.05	188	2232991	38.93	ng/ul	0.00
79) Pyrene-d10	19.35	212	2442336	38.11	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.17	264	3007106	39.36	ng/ul	0.00

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

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