

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM051021\
 Data File : BM029960.D
 Acq On : 12 May 2021 02:14
 Operator : CG/JU
 Sample : PB136110BL
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK110

Quant Time: May 12 03:27:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 12 03:18:52 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	142061	20.00	ng/ul	0.00
20) Naphthalene-d8	10.32	136	612793	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.19	164	390237	20.00	ng/ul	-0.01
64) Phenanthrene-d10	16.94	188	756032	20.00	ng/ul	0.00
79) Chrysene-d12	21.13	240	673810	20.00	ng/ul	0.00
88) Perylene-d12	23.28	264	678577	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	18654	5.24	ng/uL	0.00
4) Pyridine-d5	3.47	84	190427	21.38	ng/ul	0.00
7) Phenol-d5	6.73	99	315965	23.92	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.89	67	213618	26.08	ng/ul	0.00
11) 2-Chlorophenol-d4	7.08	132	273156	26.79	ng/ul	0.00
15) 4-Methylphenol-d8	8.26	113	262069	23.99	ng/ul	-0.01
21) Nitrobenzene-d5	8.70	128	120713	28.59	ng/ul	0.00
24) 2-Nitrophenol-d4	9.42	143	135246	28.38	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.95	165	234865	24.57	ng/ul	0.00
31) 4-Chloroaniline-d4	10.47	131	344100	23.66	ng/ul	0.00
46) Dimethylphthalate-d6	13.62	166	851432	28.15	ng/ul	0.00
49) Acenaphthylene-d8	13.88	160	1029174	26.85	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.43	143	71805	14.58	ng/ul	0.00
60) Fluorene-d10	15.19	176	718523	27.98	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.33	200	102946	20.96	ng/ul	0.00
73) Anthracene-d10	17.04	188	1101969	28.95	ng/ul	0.00
81) Pyrene-d10	19.34	212	1194525	34.56	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.14	264	1164525	30.27	ng/ul	0.00

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

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