

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052021\
 Data File : BM030117.D
 Acq On : 21 May 2021 17:48
 Operator : CG/JU
 Sample : PB136568BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK568

Quant Time: May 21 18:18:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM050721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 20 12:00:20 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	132458	20.00	ng/ul	0.00
20) Naphthalene-d8	10.28	136	575763	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.16	164	382231	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.91	188	769889	20.00	ng/ul	0.00
79) Chrysene-d12	21.10	240	659605	20.00	ng/ul	0.00
88) Perylene-d12	23.24	264	609390	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.03	96	25766	7.77	ng/uL	0.00
4) Pyridine-d5	3.45	84	241676	29.10	ng/ul	0.00
7) Phenol-d5	6.70	99	395146	32.08	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.86	67	266095	34.84	ng/ul	0.00
11) 2-Chlorophenol-d4	7.05	132	339006	35.66	ng/ul	0.00
15) 4-Methylphenol-d8	8.23	113	328918	32.29	ng/ul	0.00
21) Nitrobenzene-d5	8.67	128	157469	39.69	ng/ul	0.00
24) 2-Nitrophenol-d4	9.38	143	175050	39.09	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.92	165	290304	32.33	ng/ul	0.00
31) 4-Chloroaniline-d4	10.43	131	428646	31.37	ng/ul	0.00
46) Dimethylphthalate-d6	13.59	166	1131838	38.20	ng/ul	0.00
49) Acenaphthylene-d8	13.85	160	1361351	36.26	ng/ul	0.00
54) 4-Nitrophenol-d4	14.40	143	111071	23.02	ng/ul	0.00
60) Fluorene-d10	15.16	176	964093	38.33	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.30	200	159656	31.93	ng/ul	0.00
73) Anthracene-d10	17.01	188	1483294	38.26	ng/ul	0.00
81) Pyrene-d10	19.31	212	1579780	46.69	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.11	264	1293608	37.45	ng/ul	0.00

Target Compounds Ovalue

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

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