

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101623\
 Data File : BM042310.D
 Acq On : 16 Oct 2023 23:16
 Operator : MA/JU
 Sample : 04900-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AW3

Quant Time: Oct 17 01:30:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 15:38:21 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	114374	20.000	ng/u1	0.00
20) Naphthalene-d8	10.839	136	511104	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	306893	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	646695	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	558619	20.000	ng/u1	-0.01
88) Perylene-d12	24.062	264	612648	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	18155	4.960	ng/uL	0.00
4) Pyridine-d5	3.811	84	72290	6.668	ng/u1	0.00
7) Phenol-d5	7.163	99	99899	7.881	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	391907	45.758	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	295555	33.895	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	193285	19.255	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	199418	45.941	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	202529	43.647	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	326806	38.594	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	393359	29.503	ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	1248633	46.935	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1357653	46.628	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	34302	6.411	ng/u1	0.00
60) Fluorene-d10	15.645	176	1054005	48.196	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	164704	36.357	ng/u1	0.00
73) Anthracene-d10	17.510	188	1704090	51.440	ng/u1	0.00
81) Pyrene-d10	19.798	212	2042473	59.991	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.903	264	1899132	56.475	ng/u1	-0.01

Target Compounds	Qvalue

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

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