

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101623\
 Data File : BM042309.D
 Acq On : 16 Oct 2023 21:26
 Operator : MA/JU
 Sample : 04900-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AT6

Quant Time: Oct 17 01:30:26 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	134584	20.000	ng/u1	0.00
20) Naphthalene-d8	10.840	136	610474	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	369296	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	787467	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	690291	20.000	ng/u1	-0.01
88) Perylene-d12	24.062	264	733853	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	20923	4.858	ng/uL	0.00
4) Pyridine-d5	3.811	84	157137	12.318	ng/u1	0.00
7) Phenol-d5	7.163	99	113527	7.611	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.352	67	459935	45.637	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	341484	33.282	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	219400	18.574	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	233551	45.047	ng/u1	0.00
24) 2-Nitrophenol-d4	9.928	143	240971	43.479	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	375629	37.139	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	544001	34.160	ng/u1	0.00
46) Dimethylphthalate-d6	14.075	166	1424248	44.490	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1587841	45.319	ng/u1	0.00
54) 4-Nitrophenol-d4	14.869	143	36335	5.643	ng/u1	0.00
60) Fluorene-d10	15.645	176	1223759	46.502	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	187168	33.930	ng/u1	0.00
73) Anthracene-d10	17.510	188	1943324	48.175	ng/u1	0.00
81) Pyrene-d10	19.798	212	2266819	53.880	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.903	264	2039871	50.642	ng/u1	-0.01

Target Compounds	Qvalue

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

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