

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048340.D
 Acq On : 31 Oct 2024 02:51
 Operator : RC/JU
 Sample : P4534-17
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1L64

Quant Time: Oct 31 11:21:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM102624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Oct 27 23:39:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	272263	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.498	136	1161232	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.357	164	724958	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	1353144	20.000	ng/u1	-0.01
79) Chrysene-d12	21.321	240	1028991	20.000	ng/u1	-0.01
88) Perylene-d12	24.268	264	1076222	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	29712	4.296	ng/uL	0.00
4) Pyridine-d5	3.587	84	121264	6.050	ng/u1	0.00
7) Phenol-d5	6.892	99	168403	7.200	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.045	67	412311	28.456	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.245	132	485800	25.539	ng/u1	-0.01
15) 4-Methylphenol-d8	8.428	113	327822	17.457	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	265246	27.861	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.586	143	293192	28.427	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.127	165	497155	26.415	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.639	131	519779	19.889	ng/u1	-0.01
46) Dimethylphthalate-d6	13.768	166	1593813	29.419	ng/u1	-0.01
49) Acenaphthylene-d8	14.045	160	1752446	28.688	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.592	143	38109	3.846	ng/u1	0.01
60) Fluorene-d10	15.351	176	1346279	29.409	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.474	200	220524	24.660	ng/u1	-0.01
73) Anthracene-d10	17.198	188	2003916	31.318	ng/u1	-0.01
81) Pyrene-d10	19.486	212	2104251	33.699	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.062	264	1819601	32.055	ng/u1	-0.02

Target Compounds Qvalue

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

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