

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048342.D
 Acq On : 31 Oct 2024 04:09
 Operator : RC/JU
 Sample : P4558-02
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E1L79

Quant Time: Oct 31 11:21:48 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	282578	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.498	136	1204923	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.351	164	742118	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.098	188	1409714	20.000	ng/u1	-0.01
79) Chrysene-d12	21.321	240	1061385	20.000	ng/u1	-0.01
88) Perylene-d12	24.262	264	1132629	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	27314	3.805	ng/uL	0.00
4) Pyridine-d5	3.587	84	204107	9.811	ng/u1	0.00
7) Phenol-d5	6.893	99	166899	6.875	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	440719	29.306	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.246	132	510981	25.882	ng/u1	-0.01
15) 4-Methylphenol-d8	8.422	113	344038	17.651	ng/u1	-0.02
21) Nitrobenzene-d5	8.869	128	287632	29.116	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.587	143	313159	29.262	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	567597	29.064	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.640	131	431179	15.901	ng/u1	-0.01
46) Dimethylphthalate-d6	13.769	166	1824803	32.903	ng/u1	-0.01
49) Acenaphthylene-d8	14.045	160	2031426	32.486	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.592	143	44645	4.402	ng/u1	0.01
60) Fluorene-d10	15.351	176	1550424	33.085	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.474	200	272164	29.214	ng/u1	-0.01
73) Anthracene-d10	17.198	188	2312448	34.690	ng/u1	-0.01
81) Pyrene-d10	19.486	212	2451160	38.057	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.062	264	2118589	35.463	ng/u1	-0.02

Target Compounds Qvalue

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

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