

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047133.D
 Acq On : 09 Aug 2024 23:07
 Operator : RC/JU
 Sample : P3529-04
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AA2

Quant Time: Aug 10 00:04:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 09 01:30:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.381	152	232248	20.000	ng/u1	0.00
20) Naphthalene-d8	10.139	136	900748	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.039	164	547713	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.798	188	1088929	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.027	240	946173	20.000	ng/u1	# 0.00
88) Perylene-d12	23.733	264	1126805	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.010	96	24752	4.222	ng/uL	0.00
4) Pyridine-d5	3.410	84	95364	5.964	ng/u1	0.00
7) Phenol-d5	6.587	99	135354	7.672	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.740	67	377250	32.498	ng/u1	0.00
11) 2-Chlorophenol-d4	6.928	132	400692	26.811	ng/u1	0.00
15) 4-Methylphenol-d8	8.104	113	252282	18.051	ng/u1	0.00
21) Nitrobenzene-d5	8.528	128	219201	30.327	ng/u1	0.00
24) 2-Nitrophenol-d4	9.239	143	239019	30.281	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.775	165	414770	27.369	ng/u1	0.00
31) 4-Chloroaniline-d4	10.292	131	27042	1.358	ng/u1	0.00
46) Dimethylphthalate-d6	13.469	166	1146006	27.365	ng/u1	0.00
49) Acenaphthylene-d8	13.727	160	1323579	28.314	ng/u1	0.00
54) 4-Nitrophenol-d4	14.286	143	37204	4.595	ng/u1	0.00
60) Fluorene-d10	15.045	176	994476	27.760	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.186	200	154862	22.510	ng/u1	0.00
73) Anthracene-d10	16.898	188	1559347	30.207	ng/u1	0.00
81) Pyrene-d10	19.209	212	1712645	31.620	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.550	264	1808891	30.487	ng/u1	0.00

Target Compounds Qvalue

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

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