

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012725\
 Data File : BM049451.D
 Acq On : 27 Jan 2025 16:44
 Operator : RC/JU
 Sample : PB166231BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK231

Quant Time: Jan 28 01:49:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM011325.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 13 16:56:06 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.516	152	151523	20.000	ng/u1	0.00
20) Naphthalene-d8	10.275	136	529635	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.157	164	329963	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.904	188	672799	20.000	ng/u1	-0.01
79) Chrysene-d12	21.139	240	619555	20.000	ng/u1	0.00
88) Perylene-d12	23.915	264	668919	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.116	96	28926	7.899	ng/uL	0.00
4) Pyridine-d5	3.510	84	409925	35.161	ng/u1	0.00
7) Phenol-d5	6.710	99	414188	32.563	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.863	67	279765	33.032	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.057	132	335885	34.520	ng/u1	0.00
15) 4-Methylphenol-d8	8.228	113	298023	30.215	ng/u1	-0.01
21) Nitrobenzene-d5	8.657	128	137658	34.084	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.375	143	148941	32.988	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.910	165	293497	31.856	ng/u1	0.00
31) 4-Chloroaniline-d4	10.422	131	354432	29.236	ng/u1	0.00
46) Dimethylphthalate-d6	13.574	166	854301	34.786	ng/u1	-0.01
49) Acenaphthylene-d8	13.845	160	991077	35.350	ng/u1	0.00
54) 4-Nitrophenol-d4	14.386	143	79622	19.579	ng/u1	0.00
60) Fluorene-d10	15.157	176	754009	35.540	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.286	200	95504	22.273	ng/u1	0.00
73) Anthracene-d10	17.003	188	1179195	35.455	ng/u1	-0.01
81) Pyrene-d10	19.309	212	1373500	36.427	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.727	264	1271021	35.396	ng/u1	-0.02

Target Compounds Qvalue

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

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