

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102824\
 Data File : BM048333.D
 Acq On : 30 Oct 2024 22:15
 Operator : RC/JU
 Sample : PB164473BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK473

Quant Time: Oct 31 11:20:49 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	233518	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.498	136	965414	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.357	164	570003	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	1063868	20.000	ng/u1	-0.01
79) Chrysene-d12	21.321	240	887124	20.000	ng/u1	-0.01
88) Perylene-d12	24.262	264	979870	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.169	96	35733	6.024	ng/uL	0.00
4) Pyridine-d5	3.581	84	497116	28.915	ng/u1	0.00
7) Phenol-d5	6.887	99	566900	28.258	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.045	67	364182	29.305	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.245	132	484310	29.685	ng/u1	-0.01
15) 4-Methylphenol-d8	8.428	113	450791	27.988	ng/u1	-0.01
21) Nitrobenzene-d5	8.869	128	230267	29.092	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	252895	29.494	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.128	165	411759	26.315	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.645	131	601703	27.694	ng/u1	0.00
46) Dimethylphthalate-d6	13.768	166	1228324	28.836	ng/u1	-0.01
49) Acenaphthylene-d8	14.045	160	1411660	29.392	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.574	143	143935	18.477	ng/u1	0.00
60) Fluorene-d10	15.351	176	1038343	28.848	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	142378	20.251	ng/u1	0.00
73) Anthracene-d10	17.198	188	1558369	30.978	ng/u1	-0.01
81) Pyrene-d10	19.492	212	1687312	31.343	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.062	264	1612810	31.206	ng/u1	-0.02

Target Compounds	Qvalue

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

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