

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091124\
 Data File : BM047475.D
 Acq On : 11 Sep 2024 10:06
 Operator : RC/JU
 Sample : PB163198BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK198

Quant Time: Sep 12 03:54:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 11 00:35:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	387750	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	1715155	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	1107499	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2199548	20.000	ng/u1	0.00
79) Chrysene-d12	21.462	240	1903153	20.000	ng/u1	0.00
88) Perylene-d12	24.509	264	1922429	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	57164	5.946	ng/uL	0.00
4) Pyridine-d5	3.710	84	783578	25.510	ng/u1	0.00
7) Phenol-d5	7.022	99	887183	24.409	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.198	67	609299	26.932	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	721352	26.059	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	704215	24.650	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	365919	27.626	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	340989	27.250	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	630268	24.771	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	1004763	24.129	ng/u1	0.00
46) Dimethylphthalate-d6	13.910	166	2239320	27.509	ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	2574996	26.215	ng/u1	0.00
54) 4-Nitrophenol-d4	14.668	143	389222	24.132	ng/u1	0.00
60) Fluorene-d10	15.486	176	1951110	27.737	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	251735	21.649	ng/u1	0.00
73) Anthracene-d10	17.333	188	3073167	29.015	ng/u1	0.00
81) Pyrene-d10	19.615	212	3361634	30.307	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.303	264	2893112	28.320	ng/u1	0.00

Target Compounds	Qvalue

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

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