

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091124\
 Data File : BM047479.D
 Acq On : 11 Sep 2024 13:10
 Operator : RC/JU
 Sample : PB163275BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK275

Quant Time: Sep 12 03:59:34 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	390851	20.000	ng/u1	0.00
20) Naphthalene-d8	10.669	136	1754988	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.498	164	1105344	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2191210	20.000	ng/u1	0.00
79) Chrysene-d12	21.462	240	1827370	20.000	ng/u1	0.00
88) Perylene-d12	24.509	264	1779587	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	57479	5.931	ng/uL	0.00
4) Pyridine-d5	3.710	84	808155	26.101	ng/u1	0.00
7) Phenol-d5	7.022	99	908182	24.789	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.198	67	631425	27.688	ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	728015	26.091	ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	720169	25.008	ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	375392	27.698	ng/u1	0.00
24) 2-Nitrophenol-d4	9.745	143	350298	27.359	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	642918	24.695	ng/u1	0.00
31) 4-Chloroaniline-d4	10.792	131	1011209	23.732	ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2260105	27.819	ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	2629031	26.817	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	382034	23.733	ng/u1	0.00
60) Fluorene-d10	15.486	176	1952380	27.810	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	243770	21.044	ng/u1	0.00
73) Anthracene-d10	17.333	188	3076138	29.153	ng/u1	0.00
81) Pyrene-d10	19.615	212	3375025	31.690	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.303	264	2721459	28.778	ng/u1	0.00

Target Compounds Qvalue

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

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