

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091224\
 Data File : BM047498.D
 Acq On : 12 Sep 2024 10:50
 Operator : RC/JU
 Sample : PB163242BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK242

Quant Time: Sep 13 02:05:56 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	461955	20.000 ng/u1	0.00
20) Naphthalene-d8	10.663	136	2079396	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.492	164	1317539	20.000 ng/u1	-0.01
64) Phenanthrene-d10	17.227	188	2553224	20.000 ng/u1	-0.01
79) Chrysene-d12	21.456	240	1721485	20.000 ng/u1	-0.01
88) Perylene-d12	24.497	264	1606771	20.000 ng/u1	-0.02
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.299	96	63882	5.577 ng/uL	0.00
4) Pyridine-d5	3.710	84	965250	26.377 ng/u1	0.00
7) Phenol-d5	7.016	99	1118896	25.839 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	732053	27.160 ng/u1	0.00
11) 2-Chlorophenol-d4	7.398	132	899416	27.272 ng/u1	0.00
15) 4-Methylphenol-d8	8.563	113	914107	26.857 ng/u1	0.00
21) Nitrobenzene-d5	9.016	128	475272	29.597 ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	464523	30.620 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	865023	28.042 ng/u1	-0.01
31) 4-Chloroaniline-d4	10.786	131	1291359	25.579 ng/u1	0.00
46) Dimethylphthalate-d6	13.904	166	2818296	29.102 ng/u1	0.00
49) Acenaphthylene-d8	14.186	160	3339737	28.580 ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	441049	22.986 ng/u1	0.00
60) Fluorene-d10	15.486	176	2453181	29.315 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	251092	18.603 ng/u1	-0.01
73) Anthracene-d10	17.327	188	3674624	29.888 ng/u1	0.00
81) Pyrene-d10	19.609	212	3871206	38.585 ng/u1	0.00
92) Benzo(a)pyrene-d12	24.291	264	2551236	29.880 ng/u1	-0.02

Target Compounds Qvalue

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

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