

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100224\
 Data File : BM047792.D
 Acq On : 03 Oct 2024 01:49
 Operator : RC/JU
 Sample : PB163608BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK608

Quant Time: Oct 03 02:40:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	295327	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.598	136	1237272	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.439	164	783745	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.174	188	1610612	20.000	ng/ul	-0.01
79) Chrysene-d12	21.398	240	1502974	20.000	ng/ul	-0.01
88) Perylene-d12	24.403	264	1734558	20.000	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	38367	4.198	ng/uL	0.00
4) Pyridine-d5	3.664	84	566167	21.159	ng/ul	-0.01
7) Phenol-d5	6.969	99	621330	22.249	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	401023	20.237	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.340	132	515566	25.375	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	485995	23.407	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	250957	25.475	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.681	143	257591	27.286	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.216	165	467604	24.795	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.728	131	706481	24.232	ng/ul	-0.01
46) Dimethylphthalate-d6	13.845	166	1494510	26.302	ng/ul	-0.01
49) Acenaphthylene-d8	14.134	160	1691231	25.103	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	240515	21.260	ng/ul	0.00
60) Fluorene-d10	15.428	176	1278703	25.980	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.539	200	175189	18.588	ng/ul	-0.01
73) Anthracene-d10	17.274	188	2020066	25.507	ng/ul	-0.01
81) Pyrene-d10	19.563	212	2359222	27.658	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.198	264	2344020	26.046	ng/ul	-0.01

Target Compounds	Qvalue

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

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