

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030824\
 Data File : BM044558.D
 Acq On : 08 Mar 2024 09:41
 Operator : MA/JU
 Sample : PB159439BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK439

Quant Time: Mar 08 21:49:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 05 23:05:11 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	225220	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.498	136	1001114	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.363	164	643718	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	1327470	20.000	ng/u1	0.00
79) Chrysene-d12	21.309	240	1207964	20.000	ng/u1	0.00
88) Perylene-d12	23.568	264	1252405	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	38966	5.868	ng/uL	-0.01
4) Pyridine-d5	3.669	84	573172	29.326	ng/u1	-0.01
7) Phenol-d5	6.887	99	669944	28.933	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.069	67	445470	31.146	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.251	132	514418	30.231	ng/u1	-0.02
15) 4-Methylphenol-d8	8.422	113	525474	28.896	ng/u1	-0.02
21) Nitrobenzene-d5	8.881	128	269976	32.026	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.592	143	291991	32.130	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.122	165	480345	31.057	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.645	131	790187	30.761	ng/u1	-0.02
46) Dimethylphthalate-d6	13.786	166	1555278	30.800	ng/u1	0.00
49) Acenaphthylene-d8	14.057	160	1770874	30.665	ng/u1	0.00
54) 4-Nitrophenol-d4	14.569	143	285136	27.229	ng/u1	-0.01
60) Fluorene-d10	15.363	176	1343969	31.672	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.486	200	225399	26.158	ng/u1	0.00
73) Anthracene-d10	17.215	188	2103174	33.146	ng/u1	0.00
81) Pyrene-d10	19.515	212	2393111	32.301	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	2183602	33.171	ng/u1	0.00

Target Compounds Qvalue

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

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