

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022403.D
 Acq On : 16 Oct 2024 01:06
 Operator : RC/JU
 Sample : PB164093BL
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK093

Quant Time: Oct 16 01:31:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 14 11:01:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	96768	20.000	ng/u1	0.00
20) Naphthalene-d8	10.457	136	382786	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.304	164	314352	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.098	188	786999	20.000	ng/u1	-0.01
79) Chrysene-d12	21.533	240	843090	20.000	ng/u1	-0.02
88) Perylene-d12	24.821	264	914650	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	13283	5.334	ng/uL	0.00
4) Pyridine-d5	3.634	84	180738	26.439	ng/u1	0.00
7) Phenol-d5	6.887	99	213352	26.192	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.034	67	125083	26.127	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	161427	27.929	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	173955	26.760	ng/u1	0.00
21) Nitrobenzene-d5	8.834	128	79229	26.919	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	96134	28.212	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	192067	26.520	ng/u1	0.00
31) 4-Chloroaniline-d4	10.587	131	232796	27.203	ng/u1	0.00
46) Dimethylphthalate-d6	13.710	166	657701	26.333	ng/u1	0.00
49) Acenaphthylene-d8	13.992	160	711732	26.830	ng/u1	0.00
54) 4-Nitrophenol-d4	14.528	143	99359	23.713	ng/u1	-0.01
60) Fluorene-d10	15.304	176	598056	27.607	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.439	200	112301	21.697	ng/u1	0.00
73) Anthracene-d10	17.198	188	1041112	28.121	ng/u1	0.00
81) Pyrene-d10	19.586	212	1298436	29.123	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.598	264	1355353	27.747	ng/u1	-0.04

Target Compounds Qvalue

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

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