

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
 Data File : BP022675.D
 Acq On : 28 Oct 2024 11:37
 Operator : RC/JU
 Sample : PB164389BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK389

Quant Time: Oct 28 23:56:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 26 05:55:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	105105	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.399	136	387610	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.257	164	313171	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.063	188	769497	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.492	240	903387	20.000	ng/u1	# 0.00
88) Perylene-d12	24.745	264	1054879	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	11795	4.361	ng/uL	-0.01
4) Pyridine-d5	3.593	84	153366	20.655	ng/u1	-0.02
7) Phenol-d5	6.834	99	177718	20.087	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	6.987	67	105023	20.197	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.187	132	140020	22.304	ng/u1	-0.02
15) 4-Methylphenol-d8	8.352	113	143181	20.279	ng/u1	-0.01
21) Nitrobenzene-d5	8.787	128	66605	22.348	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.499	143	81365	23.581	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.040	165	154852	21.116	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.546	131	189130	21.826	ng/u1	-0.01
46) Dimethylphthalate-d6	13.669	166	570833	22.941	ng/u1	-0.01
49) Acenaphthylene-d8	13.946	160	571967	21.643	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.510	143	72251	17.309	ng/u1	0.00
60) Fluorene-d10	15.269	176	494558	22.915	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.410	200	99963	19.752	ng/u1	0.00
73) Anthracene-d10	17.157	188	831493	22.970	ng/u1	-0.01
81) Pyrene-d10	19.551	212	1108213	23.198	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.521	264	1340805	23.800	ng/u1	0.00

Target Compounds Qvalue

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

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