

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
 Data File : BP022503.D
 Acq On : 21 Oct 2024 13:24
 Operator : RC/JU
 Sample : PB164283BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK283

Quant Time: Oct 21 23:23:49 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 18 11:08:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	77323	20.000	ng/u1	0.00
20) Naphthalene-d8	10.422	136	284201	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.275	164	217620	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.075	188	515745	20.000	ng/u1	0.00
79) Chrysene-d12	21.504	240	588863	20.000	ng/u1	0.00
88) Perylene-d12	24.762	264	682543	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	11314	5.686	ng/uL	0.00
4) Pyridine-d5	3.611	84	145128	26.568	ng/u1	0.00
7) Phenol-d5	6.857	99	174289	26.777	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	107305	28.050	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	134030	29.020	ng/u1	0.00
15) 4-Methylphenol-d8	8.369	113	139468	26.850	ng/u1	0.00
21) Nitrobenzene-d5	8.804	128	65906	30.160	ng/u1	0.00
24) 2-Nitrophenol-d4	9.516	143	78455	31.011	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.051	165	147184	27.373	ng/u1	0.00
31) 4-Chloroaniline-d4	10.557	131	176869	27.837	ng/u1	0.00
46) Dimethylphthalate-d6	13.681	166	524548	30.338	ng/u1	0.00
49) Acenaphthylene-d8	13.963	160	531251	28.928	ng/u1	0.00
54) 4-Nitrophenol-d4	14.516	143	65935	22.731	ng/u1	0.01
60) Fluorene-d10	15.286	176	435017	29.007	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	87608	25.828	ng/u1	0.00
73) Anthracene-d10	17.169	188	739975	30.499	ng/u1	0.00
81) Pyrene-d10	19.563	212	938541	30.139	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.551	264	1116568	30.632	ng/u1	0.02

Target Compounds Qvalue

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

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