

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101724\
 Data File : BP022449.D
 Acq On : 17 Oct 2024 18:01
 Operator : RC/JU
 Sample : PB164171BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK171

Quant Time: Oct 18 04:20:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 05:42:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.664	152	181088	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	651956	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.287	164	464937	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.098	188	988215	20.000	ng/ul	0.02
79) Chrysene-d12	21.516	240	967896	20.000	ng/ul	# 0.00
88) Perylene-d12	24.780	264	1218349	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	19617	4.210	ng/uL	0.00
4) Pyridine-d5	3.611	84	252165	19.711	ng/ul	0.00
7) Phenol-d5	6.858	99	316154	20.740	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.011	67	190770	21.293	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	253475	23.435	ng/ul	0.00
15) 4-Methylphenol-d8	8.369	113	260195	21.389	ng/ul	0.00
21) Nitrobenzene-d5	8.805	128	122518	24.441	ng/ul	0.00
24) 2-Nitrophenol-d4	9.516	143	150414	25.917	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.058	165	292814	23.739	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	324982	22.297	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	845231	22.881	ng/ul	0.01
49) Acenaphthylene-d8	13.975	160	926404	23.612	ng/ul	0.01
54) 4-Nitrophenol-d4	14.510	143	105608	17.041	ng/ul	0.00
60) Fluorene-d10	15.298	176	722478	22.549	ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.434	200	129597	19.940	ng/ul	0.01
73) Anthracene-d10	17.192	188	1063715	22.881	ng/ul	0.00
81) Pyrene-d10	19.569	212	1263915	24.693	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.551	264	1268299	19.492	ng/ul	0.00

Target Compounds Qvalue

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

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