

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022350.D
 Acq On : 14 Oct 2024 11:10
 Operator : RC/JU
 Sample : PB163884BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK884

Quant Time: Oct 15 00:59:43 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	7.699	152	125399	20.000	ng/u1	0.00
20) Naphthalene-d8	10.463	136	467906	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.322	164	370195	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.116	188	943289	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1055167	20.000	ng/u1	0.00
88) Perylene-d12	24.857	264	1159294	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	13234	4.101	ng/uL	0.00
4) Pyridine-d5	3.634	84	179012	20.207	ng/u1	0.00
7) Phenol-d5	6.887	99	225687	21.381	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	130815	21.085	ng/u1	0.00
11) 2-Chlorophenol-d4	7.246	132	168485	22.495	ng/u1	0.00
15) 4-Methylphenol-d8	8.404	113	177952	21.124	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	82318	22.881	ng/u1	0.00
24) 2-Nitrophenol-d4	9.551	143	99005	23.769	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	199883	22.579	ng/u1	0.01
31) 4-Chloroaniline-d4	10.593	131	231351	22.116	ng/u1	0.00
46) Dimethylphthalate-d6	13.728	166	668563	22.730	ng/u1	0.02
49) Acenaphthylene-d8	14.010	160	719683	23.037	ng/u1	0.01
54) 4-Nitrophenol-d4	14.545	143	101124	20.494	ng/u1	0.00
60) Fluorene-d10	15.328	176	603671	23.662	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	110984	17.889	ng/u1	0.01
73) Anthracene-d10	17.216	188	1052565	23.720	ng/u1	0.01
81) Pyrene-d10	19.598	212	1366338	24.487	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.633	264	1502967	24.276	ng/u1	0.00

Target Compounds Qvalue

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

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