

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091324\
 Data File : BP021901.D
 Acq On : 13 Sep 2024 09:42
 Operator : RC/JU
 Sample : PB163242BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK242

Quant Time: Sep 14 04:42:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 12 12:48:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	156000	20.000	ng/u1	0.03
20) Naphthalene-d8	10.534	136	609392	20.000	ng/u1	# 0.03
38) Acenaphthene-d10	14.375	164	443289	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.163	188	1038459	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	1066589	20.000	ng/u1	0.00
88) Perylene-d12	24.904	264	1190364	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	19332	4.009	ng/uL	0.02
4) Pyridine-d5	3.687	84	220629	15.538	ng/u1	0.02
7) Phenol-d5	6.940	99	230693	17.307	ng/u1	0.03
9) Bis-(2-Chloroethyl)eth...	7.099	67	147778	19.088	ng/u1	0.03
11) 2-Chlorophenol-d4	7.305	132	203333	21.914	ng/u1	0.03
15) 4-Methylphenol-d8	8.457	113	243945	22.922	ng/u1	0.03
21) Nitrobenzene-d5	8.904	128	120732	31.811	ng/u1	0.03
24) 2-Nitrophenol-d4	9.616	143	124958	28.962	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.157	165	251553	26.328	ng/u1	0.03
31) 4-Chloroaniline-d4	10.657	131	333333	23.939	ng/u1	0.02
46) Dimethylphthalate-d6	13.769	166	912139	27.389	ng/u1	0.00
49) Acenaphthylene-d8	14.063	160	942370	25.796	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	133379	21.186	ng/u1	0.00
60) Fluorene-d10	15.381	176	795110	25.118	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.492	200	123199	20.737	ng/u1	0.00
73) Anthracene-d10	17.263	188	1303124	27.014	ng/u1	0.00
81) Pyrene-d10	19.633	212	1662445	28.172	ng/u1	0.01
92) Benzo(a)pyrene-d12	24.686	264	1656192	26.735	ng/u1	0.00

Target Compounds Qvalue

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

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