

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080824\
 Data File : BP021464.D
 Acq On : 09 Aug 2024 09:12
 Operator : RC/JU
 Sample : PB162595BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK595

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/09/2024
 Supervised By :mohammad ahmed 08/12/2024

Quant Time: Aug 09 23:26:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 08 18:23:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.563	152	115070	20.000	ng/u1	-0.04
20) Naphthalene-d8	10.334	136	456821	20.000	ng/u1	-0.04
38) Acenaphthene-d10	14.222	164	339999	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.051	188	747065	20.000	ng/u1	-0.02
79) Chrysene-d12	21.516	240	744218	20.000	ng/u1	0.00
88) Perylene-d12	24.780	264	853903	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	20503	8.108	ng/uL	-0.02
4) Pyridine-d5	3.552	84	241733	33.020	ng/u1	-0.03
7) Phenol-d5	6.793	99	288562	30.759	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	6.922	67	189871	33.397	ng/u1	-0.04
11) 2-Chlorophenol-d4	7.116	132	250664	32.430	ng/u1	-0.04
15) 4-Methylphenol-d8	8.305	113	236424	30.343	ng/u1	-0.04
21) Nitrobenzene-d5	8.734	128	118131m	33.293	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.452	143	135128	33.274	ng/u1	-0.04
28) 2,4-Dichlorophenol-d3	10.016	165	223776m	30.473	ng/u1	-0.03
31) 4-Chloroaniline-d4	10.516	131	331459m	34.071	ng/u1	-0.02
46) Dimethylphthalate-d6	13.646	166	838238	33.915	ng/u1	-0.01
49) Acenaphthylene-d8	13.916	160	953016	34.339	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.610	143	112215m	31.910	ng/u1	0.04
60) Fluorene-d10	15.251	176	754237	37.197	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.451	200	116289m	25.727	ng/u1	0.00
73) Anthracene-d10	17.163	188	1211377m	36.508	ng/u1	-0.01
81) Pyrene-d10	19.551	212	1491136	32.409	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.563	264	1601005	38.016	ng/u1	0.00

Target Compounds Qvalue

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

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