

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080524\
 Data File : BP021377.D
 Acq On : 05 Aug 2024 18:30
 Operator : CG/JU
 Sample : PB162219BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK219

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/06/2024
 Supervised By :mohammad ahmed 08/07/2024

Quant Time: Aug 06 01:49:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	146362	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.381	136	567218	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.263	164	358793	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.074	188	826377	20.000	ng/u1	-0.02
79) Chrysene-d12	21.527	240	882329	20.000	ng/u1	0.00
88) Perylene-d12	24.821	264	1022158	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.157	96	20010	6.221	ng/uL	-0.02
4) Pyridine-d5	3.575	84	266640	28.635	ng/u1	-0.02
7) Phenol-d5	6.845	99	328068	27.494	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	208670	28.856	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.169	132	277985	28.275	ng/u1	-0.01
15) 4-Methylphenol-d8	8.357	113	275777	27.827	ng/u1	0.00
21) Nitrobenzene-d5	8.781	128	134154	30.450	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.498	143	151240	29.993	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.063	165	240511	26.377	ng/u1	0.01
31) 4-Chloroaniline-d4	10.563	131	346317	28.670	ng/u1	0.00
46) Dimethylphthalate-d6	13.674	166	840827	32.237	ng/u1	0.00
49) Acenaphthylene-d8	13.957	160	932507	31.840	ng/u1	0.00
54) 4-Nitrophenol-d4	14.645	143	111833m	30.136	ng/u1	0.06
60) Fluorene-d10	15.286	176	738244	34.501	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.480	200	116154	23.231	ng/u1	0.02
73) Anthracene-d10	17.186	188	1229602	33.501	ng/u1	0.00
81) Pyrene-d10	19.568	212	1587790	29.107	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.592	264	1660001	32.928	ng/u1	0.00

Target Compounds Qvalue

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

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