

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022388.D
 Acq On : 15 Oct 2024 14:15
 Operator : RC/JU
 Sample : PB164092BL
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK092

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/16/2024
 Supervised By :mohammad ahmed 10/22/2024

Quant Time: Oct 16 02:16:49 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	7.646	152	87702m	20.000	ng/u1	-0.05
20) Naphthalene-d8	10.416	136	319072	20.000	ng/u1	-0.04
38) Acenaphthene-d10	14.287	164	254111	20.000	ng/u1	-0.02
64) Phenanthrene-d10	17.098	188	642836	20.000	ng/u1	-0.01
79) Chrysene-d12	21.551	240	732711	20.000	ng/u1	0.00
88) Perylene-d12	24.851	264	805691	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	10732m	4.755	ng/uL	-0.02
4) Pyridine-d5	3.599	84	157792	25.468	ng/u1	-0.04
7) Phenol-d5	6.834	99	184028	24.928	ng/u1	-0.05
9) Bis-(2-Chloroethyl)eth...	6.987	67	106207	24.477	ng/u1	-0.05
11) 2-Chlorophenol-d4	7.187	132	139584	26.646	ng/u1	-0.05
15) 4-Methylphenol-d8	8.352	113	149149m	25.316	ng/u1	-0.05
21) Nitrobenzene-d5	8.793	128	65638	26.754	ng/u1	-0.04
24) 2-Nitrophenol-d4	9.510	143	81366	28.647	ng/u1	-0.04
28) 2,4-Dichlorophenol-d3	10.046	165	161276	26.715	ng/u1	-0.04
31) 4-Chloroaniline-d4	10.551	131	188517	26.428	ng/u1	-0.04
46) Dimethylphthalate-d6	13.698	166	509444	25.233	ng/u1	-0.01
49) Acenaphthylene-d8	13.975	160	566184	26.403	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.516	143	80709	23.829	ng/u1	-0.02
60) Fluorene-d10	15.298	176	464903	26.548	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.445	200	88841	21.013	ng/u1	0.00
73) Anthracene-d10	17.198	188	829518	27.431	ng/u1	0.00
81) Pyrene-d10	19.592	212	1074153	27.722	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.627	264	1214206	28.219	ng/u1	0.00

Target Compounds Qvalue

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

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