

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023552.D
 Acq On : 20 Dec 2024 10:32
 Operator : RC/JU
 Sample : PB165707BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK707

Quant Time: Dec 21 00:31:38 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Yogesh Patel 12/21/2024
 Supervised By :mohammad ahmed 12/21/2024

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.493	152	266833m	20.000	ng/u1	-0.05
20) Naphthalene-d8	10.240	136	961255	20.000	ng/u1	#-0.04
38) Acenaphthene-d10	14.122	164	700853	20.000	ng/u1	-0.02
64) Phenanthrene-d10	16.939	188	1591571	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.380	240	1808894	20.000	ng/u1	# 0.00
88) Perylene-d12	24.545	264	1983526	20.000	ng/u1	-0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.075	96	40177	5.543	ng/uL	-0.03
4) Pyridine-d5	3.481	84	529593	28.416	ng/u1	-0.04
7) Phenol-d5	6.705	99	661700	31.605	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	6.846	67	409092	30.971	ng/u1	-0.05
11) 2-Chlorophenol-d4	7.040	132	479804	31.051	ng/u1	-0.05
15) 4-Methylphenol-d8	8.210	113	503637	29.805	ng/u1	-0.05
21) Nitrobenzene-d5	8.634	128	227938	31.755	ng/u1	-0.05
24) 2-Nitrophenol-d4	9.351	143	276397	33.410	ng/u1	-0.04
28) 2,4-Dichlorophenol-d3	9.899	165	482174	27.616	ng/u1	-0.04
31) 4-Chloroaniline-d4	10.393	131	592397	30.286	ng/u1	-0.05
46) Dimethylphthalate-d6	13.539	166	1695438	31.162	ng/u1	-0.02
49) Acenaphthylene-d8	13.810	160	1790845	31.206	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.434	143	240851	36.124	ng/u1	0.02
60) Fluorene-d10	15.145	176	1456978	30.445	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.310	200	302920	31.492	ng/u1	0.00
73) Anthracene-d10	17.039	188	2476729	34.097	ng/u1	-0.01
81) Pyrene-d10	19.439	212	3038803	32.154	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.327	264	3270034	30.796	ng/u1	-0.02

Target Compounds Qvalue

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

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