

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023500.D
 Acq On : 18 Dec 2024 16:21
 Operator : RC/JU
 Sample : P5235-01 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AS8

Manual Integrations
 APPROVED

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

Quant Time: Dec 18 16:45:24 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.528	152	259553	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.269	136	924865	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.139	164	690407	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.951	188	1554042	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.380	240	1641470	20.000	ng/u1	# 0.00
88) Perylene-d12	24.562	264	1707758	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	26671	3.783	ng/uL	-0.01
4) Pyridine-d5	3.516	84	120636	6.654	ng/u1	0.00
7) Phenol-d5	6.751	99	199687	9.805	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.881	67	441244	34.342	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.081	132	434581	28.913	ng/u1	0.00
15) 4-Methylphenol-d8	8.245	113	346909	21.106	ng/u1	-0.01
21) Nitrobenzene-d5	8.675	128	212647	30.790	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.381	143	256937	32.280	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.928	165	480699	28.614	ng/u1	0.00
31) 4-Chloroaniline-d4	10.428	131	469425	24.944	ng/u1	-0.01
46) Dimethylphthalate-d6	13.557	166	1750622	32.663	ng/u1	0.00
49) Acenaphthylene-d8	13.839	160	1547209	27.369	ng/u1	0.00
54) 4-Nitrophenol-d4	14.480	143	71929m	10.952	ng/u1	0.07
60) Fluorene-d10	15.151	176	1407286	29.852	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.322	200	302148	32.171	ng/u1	0.01
73) Anthracene-d10	17.057	188	2549314	35.944	ng/u1	0.00
81) Pyrene-d10	19.433	212	3083187	35.952	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.333	264	3169803	34.673	ng/u1	-0.01

Target Compounds Qvalue

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

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