

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023507.D
 Acq On : 18 Dec 2024 21:03
 Operator : RC/JU
 Sample : P5235-08 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT7

Manual Integrations
 APPROVED

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

Quant Time: Dec 19 04:45:19 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	221662	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.275	136	811487	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.145	164	594023	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.951	188	1422935	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.386	240	1640962	20.000	ng/u1	# 0.00
88) Perylene-d12	24.563	264	1743413	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	25462	4.229	ng/uL	-0.01
4) Pyridine-d5	3.517	84	139314	8.998	ng/u1	0.00
7) Phenol-d5	6.752	99	153458	8.823	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.881	67	352178	32.095	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.081	132	344334	26.825	ng/u1	0.00
15) 4-Methylphenol-d8	8.252	113	278992	19.875	ng/u1	0.00
21) Nitrobenzene-d5	8.675	128	182234	30.073	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.381	143	218987	31.356	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.934	165	401483	27.238	ng/u1	0.00
31) 4-Chloroaniline-d4	10.434	131	423632	25.655	ng/u1	0.00
46) Dimethylphthalate-d6	13.557	166	1436579	31.153	ng/u1	0.00
49) Acenaphthylene-d8	13.840	160	1419613	29.186	ng/u1	0.00
54) 4-Nitrophenol-d4	14.481	143	60508m	10.708	ng/u1	0.07
60) Fluorene-d10	15.151	176	1184620	29.206	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.322	200	244125	28.388	ng/u1	0.01
73) Anthracene-d10	17.057	188	2085320	32.111	ng/u1	0.00
81) Pyrene-d10	19.433	212	2724623	31.780	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.345	264	2810902	30.118	ng/u1	0.00

Target Compounds Qvalue

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

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