

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023505.D
 Acq On : 18 Dec 2024 19:43
 Operator : RC/JU
 Sample : P5235-06 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT5

Manual Integrations
 APPROVED

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

Quant Time: Dec 19 03:59:58 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	237293	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.269	136	915137	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.134	164	672550	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.945	188	1460867	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.380	240	1580730	20.000	ng/u1	# 0.00
88) Perylene-d12	24.557	264	1639821	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	30633	4.753	ng/uL	-0.01
4) Pyridine-d5	3.517	84	180833	10.911	ng/u1	0.00
7) Phenol-d5	6.752	99	178643	9.595	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.875	67	447774	38.119	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.081	132	427827	31.134	ng/u1	0.00
15) 4-Methylphenol-d8	8.246	113	332524	22.129	ng/u1	-0.01
21) Nitrobenzene-d5	8.669	128	235766	34.501	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.381	143	283085	35.943	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.928	165	514981	30.981	ng/u1	0.00
31) 4-Chloroaniline-d4	10.434	131	495640	26.616	ng/u1	0.00
46) Dimethylphthalate-d6	13.546	166	1816507	34.792	ng/u1	-0.01
49) Acenaphthylene-d8	13.822	160	1712851	31.103	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.463	143	68446m	10.698	ng/u1	0.05
60) Fluorene-d10	15.151	176	1432798	31.200	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.304	200	292242	33.101	ng/u1	0.00
73) Anthracene-d10	17.039	188	2515805	37.734	ng/u1	-0.01
81) Pyrene-d10	19.433	212	3110166	37.660	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.327	264	3161800	36.018	ng/u1	-0.02

Target Compounds Qvalue

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

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