

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023503.D
 Acq On : 18 Dec 2024 18:22
 Operator : RC/JU
 Sample : P5235-04 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT2

Quant Time: Dec 19 03:53:11 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	245460	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.263	136	874532	20.000	ng/u1	#-0.02
38) Acenaphthene-d10	14.140	164	677272	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.951	188	1491365	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.380	240	1616145	20.000	ng/u1	# 0.00
88) Perylene-d12	24.551	264	1655270	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	27005	4.050	ng/uL	-0.01
4) Pyridine-d5	3.517	84	140513	8.196	ng/u1	0.00
7) Phenol-d5	6.752	99	164774	8.556	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.875	67	356148	29.310	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.081	132	354000	24.904	ng/u1	0.00
15) 4-Methylphenol-d8	8.246	113	310925	20.003	ng/u1	-0.01
21) Nitrobenzene-d5	8.669	128	187327	28.685	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.375	143	226835	30.139	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	9.922	165	411287	25.892	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.428	131	413532	23.238	ng/u1	-0.01
46) Dimethylphthalate-d6	13.546	166	1615140	30.720	ng/u1	-0.01
49) Acenaphthylene-d8	13.828	160	1389283	25.052	ng/u1	0.00
54) 4-Nitrophenol-d4	14.481	143	72435m	11.243	ng/u1	0.07
60) Fluorene-d10	15.157	176	1286922	27.828	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.310	200	278807	30.933	ng/u1	0.00
73) Anthracene-d10	17.045	188	2436968	35.804	ng/u1	0.00
81) Pyrene-d10	19.439	212	2990966	35.423	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.339	264	2928807	33.053	ng/u1	0.00

Target Compounds Qvalue

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

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