

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023513.D
 Acq On : 19 Dec 2024 01:46
 Operator : RC/JU
 Sample : PB165677BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK677

Manual Integrations
 APPROVED

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

Quant Time: Dec 19 05:30:52 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	7.522	152	248868	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.269	136	900936	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.140	164	689011	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.945	188	1487734	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.380	240	1638258	20.000	ng/u1	# 0.00
88) Perylene-d12	24.539	264	1691904	20.000	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	35655	5.275	ng/uL	-0.01
4) Pyridine-d5	3.505	84	523296	30.105	ng/u1	-0.01
7) Phenol-d5	6.746	99	577769	29.589	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.875	67	364406	29.579	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.075	132	443965	30.805	ng/u1	-0.01
15) 4-Methylphenol-d8	8.246	113	445982	28.299	ng/u1	-0.01
21) Nitrobenzene-d5	8.675	128	186097	27.662	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.387	143	221156	28.523	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.922	165	423038	25.851	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.428	131	515757	28.133	ng/u1	-0.01
46) Dimethylphthalate-d6	13.551	166	1508598	28.204	ng/u1	0.00
49) Acenaphthylene-d8	13.828	160	1545787	27.399	ng/u1	0.00
54) 4-Nitrophenol-d4	14.440	143	208065m	31.743	ng/u1	0.03
60) Fluorene-d10	15.151	176	1139334	24.217	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.310	200	191310	21.277	ng/u1	0.00
73) Anthracene-d10	17.039	188	1949174	28.707	ng/u1	-0.01
81) Pyrene-d10	19.439	212	2405836	28.108	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.333	264	2413626	26.649	ng/u1	-0.01

Target Compounds Qvalue

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

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