

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023099.D
 Acq On : 14 Nov 2024 08:58
 Operator : RC/JU
 Sample : PB164929BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK929

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 15 00:58:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.505	152	79675m	20.000	ng/u1	-0.08
20) Naphthalene-d8	10.269	136	284671	20.000	ng/u1	-0.06
38) Acenaphthene-d10	14.175	164	220988	20.000	ng/u1	-0.03
64) Phenanthrene-d10	17.004	188	576916	20.000	ng/u1	-0.01
79) Chrysene-d12	21.439	240	702175	20.000	ng/u1	0.00
88) Perylene-d12	24.657	264	824881	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	10757	6.490	ng/uL	-0.04
4) Pyridine-d5	3.505	84	150811	30.623	ng/u1	-0.04
7) Phenol-d5	6.711	99	172731	29.323	ng/u1	-0.07
9) Bis-(2-Chloroethyl)eth...	6.858	67	111253	32.146	ng/u1	-0.07
11) 2-Chlorophenol-d4	7.058	132	134955	31.282	ng/u1	-0.07
15) 4-Methylphenol-d8	8.216	113	140081m	28.898	ng/u1	-0.08
21) Nitrobenzene-d5	8.652	128	65133	32.865	ng/u1	-0.07
24) 2-Nitrophenol-d4	9.369	143	81300	34.410	ng/u1	-0.06
28) 2,4-Dichlorophenol-d3	9.916	165	152966	30.916	ng/u1	-0.05
31) 4-Chloroaniline-d4	10.416	131	171115	28.989	ng/u1	-0.06
46) Dimethylphthalate-d6	13.575	166	542557	34.065	ng/u1	-0.04
49) Acenaphthylene-d8	13.857	160	564686	33.978	ng/u1	-0.04
54) 4-Nitrophenol-d4	14.440	143	76942m	32.145	ng/u1	-0.01
60) Fluorene-d10	15.204	176	473285	33.893	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.345	200	97966	28.347	ng/u1	-0.02
73) Anthracene-d10	17.104	188	843074	34.514	ng/u1	-0.01
81) Pyrene-d10	19.486	212	1128833	34.902	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.427	264	1368362	34.825	ng/u1	-0.02

Target Compounds Qvalue

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

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