

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014264.D
 Acq On : 23 Mar 2023 19:17
 Operator : CG/JU
 Sample : PB151603BL
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK603

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/24/2023
 Supervised By : Jagrut Upadhyay 03/24/2023

Quant Time: Mar 24 00:53:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 23 04:56:06 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	155341	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	686280	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.681	164	511137	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.439	188	1190108	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.521	240	1150781	20.000	ng/u1	# 0.00
88) Perylene-d12	24.039	264	1059219	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	25080	6.100	ng/uL	0.00
4) Pyridine-d5	3.834	84	347009	30.483	ng/u1	0.00
7) Phenol-d5	7.193	99	451013	32.900	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.357	67	291777	36.496	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	337335	32.299	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	374883	33.479	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	177184	32.358	ng/u1	0.00
24) 2-Nitrophenol-d4	9.934	143	208188	31.761	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	359360	29.137	ng/u1	0.00
31) 4-Chloroaniline-d4	10.998	131	525492	32.196	ng/u1	0.00
46) Dimethylphthalate-d6	14.087	166	1388009	31.997	ng/u1	0.00
49) Acenaphthylene-d8	14.375	160	1455255	31.668	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	200336m	26.207	ng/u1	0.00
60) Fluorene-d10	15.675	176	1168697	31.789	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	228657	25.421	ng/u1	0.00
73) Anthracene-d10	17.539	188	1959706	33.920	ng/u1	0.00
81) Pyrene-d10	19.763	212	2356062	37.121	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.874	264	2015649	35.775	ng/u1	0.00

Target Compounds Qvalue

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

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