

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032423\
 Data File : BP014282.D
 Acq On : 24 Mar 2023 09:45
 Operator : CG/JU
 Sample : PB151650BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK650

Quant Time: Mar 24 22:33:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 24 22:30:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	163021	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	733193	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.681	164	530073	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.439	188	1245012	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.522	240	1156009	20.000	ng/u1	#-0.01
88) Perylene-d12	24.039	264	1014541	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	23919	5.544	ng/uL	0.00
4) Pyridine-d5	3.834	84	331265	27.729	ng/u1	0.00
7) Phenol-d5	7.193	99	435412	30.266	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	283130	33.746	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	324566	29.612	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	356964	30.377	ng/u1	-0.01
21) Nitrobenzene-d5	9.210	128	168007	28.719	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	203110	29.004	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	338110	25.660	ng/u1	0.00
31) 4-Chloroaniline-d4	11.004	131	487396	27.951	ng/u1	0.00
46) Dimethylphthalate-d6	14.087	166	1338403	29.752	ng/u1	0.00
49) Acenaphthylene-d8	14.375	160	1414628	29.685	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	184607	23.287	ng/u1	0.00
60) Fluorene-d10	15.675	176	1130937	29.663	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	212998	22.636	ng/u1	0.00
73) Anthracene-d10	17.539	188	1867591	30.900	ng/u1	0.00
81) Pyrene-d10	19.763	212	2252656	35.332	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.880	264	1791320	33.194	ng/u1	-0.01

Target Compounds Qvalue

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

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