

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032023\
 Data File : BP014219.D
 Acq On : 22 Mar 2023 14:22
 Operator : CG/JU
 Sample : PB151568BL
 Misc :
 ALS Vial : 91 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK568

Quant Time: Mar 23 00:59:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 22 03:05:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	187725	20.000	ng/u1	0.00
20) Naphthalene-d8	10.863	136	806753	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.681	164	560744	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.434	188	1288920	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.516	240	997361	20.000	ng/u1	# 0.00
88) Perylene-d12	24.033	264	910105	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	26839	5.402	ng/uL	0.00
4) Pyridine-d5	3.846	84	370284	26.917	ng/u1	0.00
7) Phenol-d5	7.205	99	500669	30.222	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	307243	31.801	ng/u1	0.00
11) 2-Chlorophenol-d4	7.569	132	377415	29.903	ng/u1	0.00
15) 4-Methylphenol-d8	8.758	113	409012	30.226	ng/u1	0.00
21) Nitrobenzene-d5	9.216	128	192015	29.830	ng/u1	0.00
24) 2-Nitrophenol-d4	9.940	143	227320	29.501	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	400158	27.600	ng/u1	0.00
31) 4-Chloroaniline-d4	11.004	131	538452	28.063	ng/u1	0.00
46) Dimethylphthalate-d6	14.087	166	1378940	28.976	ng/u1	0.00
49) Acenaphthylene-d8	14.375	160	1488008	29.517	ng/u1	0.00
54) 4-Nitrophenol-d4	14.892	143	174505	20.808	ng/u1	0.00
60) Fluorene-d10	15.675	176	1178926	29.231	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	214205	21.988	ng/u1	0.00
73) Anthracene-d10	17.534	188	1932067	30.878	ng/u1	0.00
81) Pyrene-d10	19.763	212	2196456	39.930	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.868	264	1539401	31.799	ng/u1	0.00

Target Compounds Qvalue

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

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