

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100322\
 Data File : BP011886.D
 Acq On : 03 Oct 2022 23:38
 Operator : CG/JU
 Sample : PB148023BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK023

Quant Time: Oct 04 00:50:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 01 00:53:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.887	152	253858	20.000 ng/u1	0.00
20) Naphthalene-d8	10.699	136	1015988	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.534	164	585455	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.292	188	1198564	20.000 ng/u1	0.00
79) Chrysene-d12	21.374	240	1292355	20.000 ng/u1	-0.01
88) Perylene-d12	23.810	264	1359096	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.305	96	35826	6.224 ng/uL	0.00
4) Pyridine-d5	3.722	84	397044	23.549 ng/u1	-0.01
7) Phenol-d5	7.052	99	588133	26.836 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	339660	27.961 ng/u1	-0.01
11) 2-Chlorophenol-d4	7.416	132	509106	29.428 ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	463812	25.602 ng/u1	0.00
21) Nitrobenzene-d5	9.057	128	235318	30.408 ng/u1	0.00
24) 2-Nitrophenol-d4	9.775	143	244549	29.537 ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.316	165	436265	28.487 ng/u1	0.00
31) 4-Chloroaniline-d4	10.840	131	627515	26.878 ng/u1	0.00
46) Dimethylphthalate-d6	13.945	166	1284151	30.673 ng/u1	0.00
49) Acenaphthylene-d8	14.228	160	1518137	31.808 ng/u1	0.00
54) 4-Nitrophenol-d4	14.734	143	215536	24.861 ng/u1	0.00
60) Fluorene-d10	15.528	176	1128669	31.043 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	170452	23.496 ng/u1	0.00
73) Anthracene-d10	17.386	188	1776884	32.545 ng/u1	-0.01
81) Pyrene-d10	19.622	212	2234129	34.135 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.651	264	2338720	34.375 ng/u1	-0.01

Target Compounds Qvalue

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

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