

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070722\
 Data File : BP010934.D
 Acq On : 07 Jul 2022 11:42
 Operator : CG/JU
 Sample : PB146060BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK060

Quant Time: Jul 08 00:58:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 06 23:13:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	336669	20.000	ng/u1	0.00
20) Naphthalene-d8	10.498	136	1300517	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.363	164	670467	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.128	188	1306184	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.239	240	1294968	20.000	ng/u1	# 0.00
88) Perylene-d12	23.598	264	1324432	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	67834	7.622	ng/uL	0.00
4) Pyridine-d5	3.611	84	852013	35.964	ng/u1	0.00
7) Phenol-d5	6.881	99	942744	35.374	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.046	67	617130	35.407	ng/u1	0.00
11) 2-Chlorophenol-d4	7.240	132	796622	36.028	ng/u1	0.00
15) 4-Methylphenol-d8	8.422	113	728160	34.563	ng/u1	0.00
21) Nitrobenzene-d5	8.869	128	340602	41.140	ng/u1	0.00
24) 2-Nitrophenol-d4	9.587	143	316218	46.167	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.122	165	592635	35.017	ng/u1	0.00
31) 4-Chloroaniline-d4	10.646	131	979199	35.445	ng/u1	0.00
46) Dimethylphthalate-d6	13.781	166	1666200	36.279	ng/u1	0.00
49) Acenaphthylene-d8	14.051	160	2017691	34.422	ng/u1	0.00
54) 4-Nitrophenol-d4	14.575	143	241671	33.689	ng/u1	0.00
60) Fluorene-d10	15.363	176	1459071	37.020	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	141359	32.665	ng/u1	0.00
73) Anthracene-d10	17.228	188	2218328	38.172	ng/u1	0.00
81) Pyrene-d10	19.474	212	2485200	35.226	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.445	264	2515891	38.441	ng/u1	0.00

Target Compounds Qvalue

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

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