

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011122.D
 Acq On : 27 Jul 2022 10:45
 Operator : CG/JU
 Sample : PB146439BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK439

Quant Time: Jul 28 00:00:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 26 14:31:36 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	290137	20.000	ng/u1	0.00
20) Naphthalene-d8	10.451	136	1209143	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	624124	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.092	188	1078835	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.210	240	971317	20.000	ng/u1	# 0.00
88) Perylene-d12	23.551	264	1066154	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	50683	6.298	ng/uL	0.00
4) Pyridine-d5	3.569	84	659179	31.688	ng/u1	0.00
7) Phenol-d5	6.840	99	754955	31.699	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	562041	36.699	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	628541	32.661	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	598006	31.364	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	314323	35.671	ng/u1	0.00
24) 2-Nitrophenol-d4	9.546	143	293778	33.693	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	478407	30.003	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	788620	30.000	ng/u1	0.00
46) Dimethylphthalate-d6	13.745	166	1417001	33.219	ng/u1	0.00
49) Acenaphthylene-d8	14.016	160	1824891	36.259	ng/u1	0.00
54) 4-Nitrophenol-d4	14.539	143	172364	19.965	ng/u1	0.00
60) Fluorene-d10	15.328	176	1177092	32.464	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	135878	25.453	ng/u1	0.00
73) Anthracene-d10	17.192	188	1639976	35.757	ng/u1	0.00
81) Pyrene-d10	19.445	212	1762062	34.541	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.398	264	1869094	36.501	ng/u1	0.00

Target Compounds Qvalue

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

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