

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072022\
 Data File : BP011050.D
 Acq On : 20 Jul 2022 09:27
 Operator : CG/JU
 Sample : PB146316BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK316

Quant Time: Jul 21 00:42:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	352963	20.000	ng/u1	0.00
20) Naphthalene-d8	10.469	136	1453184	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.340	164	826461	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	1667404	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.227	240	1585148	20.000	ng/u1	# 0.00
88) Perylene-d12	23.580	264	1533563	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	64666	6.605	ng/uL	0.00
4) Pyridine-d5	3.587	84	806150	31.856	ng/u1	0.00
7) Phenol-d5	6.852	99	934486	32.253	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.022	67	619346	33.243	ng/u1	0.00
11) 2-Chlorophenol-d4	7.210	132	775430	33.122	ng/u1	0.00
15) 4-Methylphenol-d8	8.393	113	739039	31.862	ng/u1	0.00
21) Nitrobenzene-d5	8.840	128	363640	34.338	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	350694	33.467	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	607638	31.708	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	977260	30.933	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	2001147	35.428	ng/u1	0.00
49) Acenaphthylene-d8	14.028	160	2410059	36.162	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	283175	24.770	ng/u1	0.00
60) Fluorene-d10	15.339	176	1674381	34.873	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	211495	25.633	ng/u1	0.00
73) Anthracene-d10	17.204	188	2536218	35.779	ng/u1	0.00
81) Pyrene-d10	19.463	212	2903271	34.873	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.427	264	2696272	36.606	ng/u1	0.00

Target Compounds Qvalue

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

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