

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072522\
 Data File : BP011083.D
 Acq On : 25 Jul 2022 09:50
 Operator : CG/JU
 Sample : PB146370BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK370

Quant Time: Jul 25 11:42:36 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	7.669	152	409816	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.463	136	1910072	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.334	164	1349746	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.104	188	2773855	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.222	240	2178272	20.000	ng/u1	# 0.00
88) Perylene-d12	23.568	264	1751254	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	78813	6.933	ng/uL	-0.02
4) Pyridine-d5	3.575	84	1090223	37.105	ng/u1	-0.02
7) Phenol-d5	6.852	99	1281995	38.109	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	875092	40.454	ng/u1	0.00
11) 2-Chlorophenol-d4	7.205	132	1058537	38.942	ng/u1	0.00
15) 4-Methylphenol-d8	8.387	113	1093733	40.612	ng/u1	-0.01
21) Nitrobenzene-d5	8.834	128	564019	40.520	ng/u1	0.00
24) 2-Nitrophenol-d4	9.557	143	592160	42.993	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	941122	37.363	ng/u1	0.00
31) 4-Chloroaniline-d4	10.616	131	1587619	38.232	ng/u1	0.00
46) Dimethylphthalate-d6	13.757	166	3872219	41.975	ng/u1	0.00
49) Acenaphthylene-d8	14.028	160	4507994	41.417	ng/u1	0.00
54) 4-Nitrophenol-d4	14.557	143	607968	32.563	ng/u1	0.00
60) Fluorene-d10	15.339	176	3150399	40.177	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.469	200	522148	38.041	ng/u1	0.00
73) Anthracene-d10	17.204	188	4745456	40.242	ng/u1	0.00
81) Pyrene-d10	19.457	212	5114597	44.707	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.421	264	3430301	40.783	ng/u1	0.00

Target Compounds Qvalue

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

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