

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072722\
 Data File : BP011145.D
 Acq On : 28 Jul 2022 00:36
 Operator : CG/JU
 Sample : PB146457BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK457

Quant Time: Jul 28 01:16:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 28 00:46:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.657	152	327527	20.000	ng/u1	0.00
20) Naphthalene-d8	10.446	136	1328012	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.322	164	658697	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.086	188	1192033	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.204	240	1192775	20.000	ng/u1	# 0.00
88) Perylene-d12	23.539	264	1348823	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	54584	6.008	ng/uL	0.00
4) Pyridine-d5	3.569	84	734313	31.271	ng/u1	0.00
7) Phenol-d5	6.834	99	951441	35.389	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.999	67	676820	39.149	ng/u1	0.00
11) 2-Chlorophenol-d4	7.193	132	781639	35.980	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	740221	34.391	ng/u1	0.00
21) Nitrobenzene-d5	8.822	128	377683	39.025	ng/u1	0.00
24) 2-Nitrophenol-d4	9.540	143	360016	37.594	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.075	165	591020	33.748	ng/u1	0.00
31) 4-Chloroaniline-d4	10.598	131	941709	32.617	ng/u1	0.00
46) Dimethylphthalate-d6	13.740	166	1645988	36.562	ng/u1	0.00
49) Acenaphthylene-d8	14.010	160	2128316	40.068	ng/u1	0.00
54) 4-Nitrophenol-d4	14.539	143	230468	25.295	ng/u1	0.00
60) Fluorene-d10	15.322	176	1364809	35.666	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.445	200	164711	27.924	ng/u1	0.00
73) Anthracene-d10	17.186	188	1987662	39.223	ng/u1	0.00
81) Pyrene-d10	19.439	212	2296667	36.662	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.392	264	2554794	39.436	ng/u1	0.00

Target Compounds Qvalue

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

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